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(Acts adopted under the EC Treaty/Euratom Treaty whose publication is obligatory)

REGULATIONS

REGULATION (EC) No 1272/2008 OF THE EUROPEAN PARLIAMENT AND OF THE COUNCIL

of 16 December 2008

on classification, labelling and packaging of substances and mixtures, amending and repealing Directives 67/548/EEC and 1999/45/EC, and amending Regulation (EC) No 1907/2006

(Text with EEA relevance)

THE EUROPEAN PARLIAMENT AND THE COUNCIL OF THE EUROPEAN UNION,

Having regard to the Treaty establishing the European Community, and in particular Article 95 thereof,

Having regard to the proposal from the Commission,

Having regard to the opinion of the European Economic and Social Committee [1],

Acting in accordance with the procedure laid down in Article 251 of the Treaty [2],

Whereas:

(1) This Regulation should ensure a high level of protection of human health and the environment as well as the free movement of chemical substances, mixtures and certain specific articles, while enhancing competitiveness and innovation.

(2) The efficient functioning of the internal market for substances, mixtures and those articles can be achieved only if the requirements applicable to them do not differ significantly between Member States.

(3) A high level of human health and environmental protection should be ensured in the approximation of legislation on the criteria for classification and labelling of substances and mixtures, with the goal of achieving sustainable development.

(4) Trade in substances and mixtures is an issue relating not only to the internal market, but also to the global market. Enterprises should therefore benefit from the global harmonisation of rules for classification and labelling and from consistency between, on the one hand, the rules for classification and labelling for supply and use and, on the other hand, those for transport.

(5) With a view to facilitating worldwide trade while protecting human health and the environment, harmonised criteria for classification and labelling have been carefully developed over a period of 12 years within the United Nations (UN) structure, resulting in the Globally Harmonised System of Classification and Labelling of Chemicals (hereinafter referred to as 'the GHS').

(6) This Regulation follows various declarations whereby the Community confirmed its intention to contribute to the global harmonisation of criteria for classification and labelling, not only at UN level, but also through the incorporation of the internationally agreed GHS criteria into Community law.

(7) The benefits for enterprises will increase as more countries in the world adopt the GHS criteria in their legislation. The Community should be at the forefront of this process to encourage other countries to follow and with the aim of providing a competitive advantage to industry in the Community.

(8) Therefore it is essential to harmonise the provisions and criteria for the classification and labelling of substances, mixtures and certain specific articles within the Community, taking into account the classification criteria and labelling rules of the GHS, but also by building on the 40 years of experience obtained through implementation of existing Community chemicals legislation and maintaining

[1] OJ C 204, 9.8.2008, p. 47.

[2] Opinion of the European Parliament of 3 September 2008 (not yet published in the Official Journal).

the level of protection achieved through the system of harmonisation of classification and labelling, through Community hazard classes not yet part of the GHS as well as through current labelling and packaging rules.

(9) This Regulation should be without prejudice to the full and complete application of Community competition rules.

(10) The objective of this Regulation should be to determine which properties of substances and mixtures should lead to a classification as hazardous, in order for the hazards of substances and mixtures to be properly identified and communicated. Such properties should include physical hazards as well as hazards to human health and to the environment, including hazards to the ozone layer.

(11) This Regulation should, as a general principle, apply to all substances and mixtures supplied in the Community, except where other Community legislation lays down more specific rules on classification and labelling, such as Council Directive 76/768/EEC of 27 July 1976 on the approximation of the laws of the Member States relating to cosmetic products ([1]), Council Directive 82/471/EEC of 30 June 1982 concerning certain products used in animal nutrition ([2]), Council Directive 88/388/EEC of 22 June 1988 on the approximation of the laws of the Member States relating to flavourings for use in foodstuffs and to source materials for their production ([3]), Council Directive 89/107/EEC of 21 December 1988 on the approximation of the laws of the Member States concerning food additives authorised for use in foodstuffs intended for human consumption ([4]), Council Directive 90/385/EEC of 20 June 1990 on the approximation of the laws of the Member States relating to active implantable medical devices ([5]), Council Directive 93/42/EEC of 14 June 1993 concerning medical devices ([6]), Directive 98/79/EC of the European Parliament and of the Council of 27 October 1998 on in vitro diagnostic medical devices ([7]), Commission Decision 1999/217/EC of 23 February 1999 adopting a register of flavouring substances used in or on foodstuffs drawn up in application of Regulation (EC) No 2232/96 of the European Parliament and of the Council ([8]), Directive 2001/82/EC of the European Parliament and of the Council of 6 November 2001 on the Community code relating to veterinary medicinal products ([9]), Directive 2001/83/EC of the European Parliament and of the Council of 6 November 2001 on the Community code relating to medicinal products for human use ([10]), Regulation (EC) No 178/2002 of the European Parliament and of the Council of 28 January 2002 laying down the general principles and requirements of food law, establishing the European Food Safety Authority and laying down procedures in matters of food safety ([11]) and Regulation (EC) No 1831/2003 of the European Parliament and of the Council of 22 September 2003 on additives for use in animal nutrition ([12]) or except where substances and mixtures are transported by air, sea, road, rail or inland waterways.

(12) The terms and definitions used in this Regulation should be consistent with those set out in Regulation (EC) No 1907/2006 of the European Parliament and of the Council of 18 December 2006 concerning the Registration, Evaluation, Authorisation and Restriction of Chemicals (REACH) ([13]), with those set out in the rules governing transport and with the definitions specified at UN level in the GHS, in order to ensure maximum consistency in the application of chemicals legislation within the Community in the context of global trade. The hazard classes specified in the GHS should be set out in this Regulation for the same reason.

(13) It is especially appropriate to include those hazard classes defined in the GHS which specifically take account of the fact that the physical hazards which may be exhibited by substances and mixtures are to some extent influenced by the way in which they are released.

(14) The term 'mixture' as defined in this Regulation should have the same meaning as the term 'preparation' previously used in Community legislation.

(15) This Regulation should replace Council Directive 67/548/EEC of 27 June 1967 on the approximation of the laws, regulations and administrative provisions relating to the classification, packaging and labelling of dangerous substances ([14]) as well as Directive 1999/45/EC of the European Parliament and of the Council of 31 May 1999 concerning the approximation of the laws, regulations and administrative provisions of the Member States relating to the classification, packaging and labelling of dangerous preparations ([15]). It should maintain the overall current level of protection of human health and the environment provided by those Directives. Therefore, some hazard classes which are covered by those Directives but are not yet included in the GHS should be maintained in this Regulation.

(16) Responsibility for the identification of hazards of substances and mixtures and for deciding on their classification should mainly lie with manufacturers, importers and downstream users of those substances or mixtures, regardless of whether they are subject to the requirements

([1]) OJ L 262, 27.9.1976, p. 169.
([2]) OJ L 213, 21.7.1982, p. 8.
([3]) OJ L 184, 15.7.1988, p. 61.
([4]) OJ L 40, 11.2.1989, p. 27.
([5]) OJ L 189, 20.7.1990, p. 17.
([6]) OJ L 169, 12.7.1993, p. 1.
([7]) OJ L 331, 7.12.1998, p. 1.
([8]) OJ L 84, 27.3.1999, p. 1.
([9]) OJ L 311, 28.11.2001, p. 1.
([10]) OJ L 311, 28.11.2001, p. 67.
([11]) OJ L 31, 1.2.2002, p. 1.
([12]) OJ L 268, 18.10.2003, p. 29.
([13]) OJ L 396, 30.12.2006, p. 1. Corrected version in OJ L 136, 29.5.2007, p. 3.
([14]) OJ 196, 16.8.1967, p. 1.
([15]) OJ L 200, 30.7.1999, p. 1.

of Regulation (EC) No 1907/2006. In fulfilling their responsibilities for classification, downstream users should be allowed to use the classification of a substance or mixture derived in accordance with this Regulation by an actor in the supply chain, provided that they do not change the composition of the substance or mixture. Responsibility for classification of substances not placed on the market that are subject to registration or notification under Regulation (EC) No 1907/2006 should mainly lie with the manufacturers, producers of articles and importers. However, there should be a possibility to provide for harmonised classifications of substances for hazard classes of highest concern and of other substances on a case-by-case basis which should be applied by all manufacturers, importers and downstream users of such substances and of mixtures containing such substances.

(17) Where a decision has been taken to harmonise the classification of a substance for a specific hazard class or differentiation within a hazard class by including or revising an entry for that purpose in Part 3 of Annex VI to this Regulation, the manufacturer, importer and downstream user should apply this harmonised classification, and only self-classify for the remaining, non-harmonised hazard classes or differentiations within the hazard class.

(18) To ensure that customers receive information on hazards, suppliers of substances and mixtures should ensure that they are labelled and packaged in accordance with this Regulation before placing them on the market, according to the classification derived. In fulfilling their responsibilities downstream users should be allowed to use the classification of a substance or mixture derived in accordance with this Regulation by an actor in the supply chain, provided that they do not change the composition of the substance or mixture, and distributors should be allowed to use the classification of a substance or mixture derived in accordance with this Regulation by an actor in the supply chain.

(19) To ensure information on hazardous substances is available when they are included in mixtures containing at least one substance that is classified as hazardous, supplemental labelling information should be provided, where applicable.

(20) While a manufacturer, importer or downstream user of any substance or mixture should not be obliged to generate new toxicological or eco-toxicological data for the purpose of classification, he should identify all relevant information available to him on the hazards of the substance or mixture and evaluate its quality. The manufacturer, importer or downstream user should also take into account historical human data, such as epidemiological studies on exposed populations, accidental or occupational exposure and effect data, and clinical studies. That information should be compared with the criteria for the different hazard classes and differentiations in order for that manufacturer, importer or downstream user to arrive at a conclusion as to whether or not the substance or mixture should be classified as hazardous.

(21) While the classification of any substance or mixture may be carried out on the basis of available information, the available information to be used for the purposes of this Regulation should preferably have been generated in accordance with the test methods referred to in Regulation (EC) No 1907/2006, transport provisions or international principles or procedures for the validation of information, so as to ensure quality and comparability of the results and consistency with other requirements at international or Community level. The same test methods, provisions, principles and procedures should be followed where the manufacturer, importer or downstream user chooses to generate new information.

(22) To facilitate hazard identification for mixtures, manufacturers, importers and downstream users should base this identification on the data for the mixture itself, where available, except for mixtures with carcinogenic, germ cell mutagenic or reproductive toxic substances, or where the biodegradation or bioaccumulation properties in the hazard class hazardous to the aquatic environment are evaluated. In those cases, as the hazards of the mixture cannot be sufficiently assessed in a manner that is based on the mixture itself, the data for the individual substances of the mixture should normally be used as a basis for the hazard identification of the mixture.

(23) If sufficient information is available on similar tested mixtures, including relevant ingredients of the mixtures, it is possible to determine the hazardous properties of an untested mixture by applying certain rules known as 'bridging principles'. Those rules allow characterisation of the hazards of the mixture without performing tests on it, but rather by building on the available information on similar tested mixtures. Where no or inadequate test data are available for the mixture itself, manufacturers, importers and downstream users should therefore follow the bridging principles to ensure adequate comparability of results of the classification of such mixtures.

(24) Specific industry sectors may establish networks to facilitate exchange of data and bring together expertise in the evaluation of information, test data, weight of evidence determinations and bridging principles. Such networks may support manufacturers, importers and downstream users within those industry sectors, and in particular small and medium-sized enterprises (SMEs) in the fulfilment of their obligations under this Regulation. Those networks may also be used to exchange information and best practices with a view to simplifying fulfilment of the notification obligations. Suppliers making use of such support should remain fully responsible for the fulfilment of their classification, labelling and packaging responsibilities under this Regulation.

(25) The protection of animals falling within the scope of Council Directive 86/609/EEC of 24 November 1986 on the approximation of laws, regulations and administrative provisions of the Member States regarding the protection of animals used for experimental and other scientific purposes [1] is of high priority. Accordingly, where the manufacturer, importer or downstream user chooses to generate information for the purposes of this Regulation, they should first consider means other than testing on animals within the scope of Directive 86/609/EEC. Tests on non-human primates should be prohibited for the purposes of this Regulation.

(26) The test methods in Commission Regulation (EC) No 440/2008 of 30 May 2008 laying down test methods pursuant to Regulation (EC) No 1907/2006 of the European Parliament and of the Council on the Registration, Evaluation, Authorisation and Restriction of Chemicals (REACH) [2] are regularly reviewed and improved with a view to reducing testing on vertebrate animals and the number of animals involved. The European Centre for the Validation of Alternative Methods (ECVAM) of the Commission's Joint Research Centre plays an important role in the scientific assessment and validation of alternative test methods.

(27) The classification and labelling criteria set out in this Regulation should take the utmost account of promoting alternative methods for the assessment of hazards of substances and mixtures and of the obligation to generate information on intrinsic properties by means other than tests on animals within the meaning of Directive 86/609/EEC as laid down in Regulation (EC) No 1907/2006. Future criteria should not become a barrier to this aim and the corresponding obligations under that Regulation, and should under no circumstances lead to the use of animal tests where alternative tests are adequate for the purposes of classification and labelling.

(28) For the purposes of classification, data should not be generated by means of testing on humans. Available, reliable epidemiological data and experience with regard to the effects of substances and mixtures on humans (e.g. occupational data and data from accident databases) should be taken into account and may be given priority over data derived from animal studies when they demonstrate hazards not identified from those studies. The results of animal studies should be weighed against the results of data from humans and expert judgement should be used to ensure the best protection of human health when evaluating both the animal and human data.

(29) New information as regards physical hazards should always be necessary, except if the data are already available or if a derogation is provided for in this Regulation.

(30) Testing that is carried out for the sole purpose of this Regulation should be carried out on the substance or mixture in the form(s) or physical state(s) in which the substance or mixture is placed on the market and in which it can reasonably be expected to be used. It should, however, be possible to use, for the purpose of this Regulation, the results of tests that are carried out to comply with other regulatory requirements, including those laid down by third countries, even if the tests were not carried out on the substance or mixture in the form(s) or physical state(s) in which it is placed on the market and in which it can reasonably be expected to be used.

(31) If tests are performed, they should comply where appropriate with the relevant requirements for the protection of laboratory animals, set out in Directive 86/609/EEC, and, in the case of ecotoxicological and toxicological tests, good laboratory practice, set out in Directive 2004/10/EC of the European Parliament and of the Council of 11 February 2004 on the harmonisation of laws, regulations and administrative provisions relating to the application of the principles of good laboratory practice and the verification of their application for tests on chemical substances [3].

(32) The criteria for classification in different hazard classes and differentiations should be set out in an annex, which should also contain additional provisions as to how the criteria may be met.

(33) Recognising that the application of the criteria for the different hazard classes to information is not always straightforward and simple, manufacturers, importers and downstream users should apply weight of evidence determinations involving expert judgement to arrive at adequate results.

(34) Specific concentration limits for substances should be assigned to a substance by a manufacturer, importer or downstream user in accordance with the criteria referred to in this Regulation, provided the manufacturer, importer or downstream user is able to justify the limits and informs the European Chemicals Agency (hereinafter referred to as 'the Agency') accordingly. However, specific concentration limits should not be set for harmonised hazard classes or differentiations for substances included in the harmonised classification and labelling tables annexed to this Regulation. Guidance should be provided by the Agency for the purpose of setting the specific concentration limits. In order to ensure uniformity, specific concentration limits should also be included, where appropriate, in cases of harmonised classifications. Specific concentration limits should take precedence over any other concentration limit for the purpose of classification.

(35) Multiplying factors (M-factors) for substances classified as hazardous to the aquatic environment, acute category 1 or chronic category 1, should be assigned to a substance by a manufacturer, importer or downstream user in accordance with the criteria referred to in this Regulation. Guidance should be provided by the Agency for the purpose of setting the M-factors.

[1] OJ L 358, 18.12.1986, p. 1.
[2] OJ L 142, 31.5.2008, p. 1.
[3] OJ L 50, 20.2.2004, p. 44.

(36) For reasons of proportionality and workability, generic cut-off values should be defined, both for identified impurities, additives and individual constituents of substances and for substances in mixtures, specifying when information on these should be taken into account in determining the hazard classification of substances and mixtures.

(37) To ensure adequate classification of mixtures, available information on synergistic and antagonistic effects should be taken into account for the classification of mixtures.

(38) Manufacturers, importers and downstream users should re-evaluate the classifications of substances or mixtures they place on the market if they become aware of new adequate and reliable scientific or technical information that may affect those classifications or if they change the composition of their mixtures, to ensure that the classification is based on up-to-date information, unless there is sufficient evidence that the classification would not change. Suppliers should update the labels accordingly.

(39) Substances and mixtures classified as hazardous should be labelled and packaged according to their classification, so as to ensure appropriate protection and to provide essential information to their recipients, by drawing their attention to the hazards of the substance or mixture.

(40) The two instruments foreseen by this Regulation to be used to communicate the hazards of substances and mixtures are labels and the safety data sheets provided for in Regulation (EC) No 1907/2006. Of these two, the label is the only tool for communication to consumers, but it may also serve to draw the attention of workers to the more comprehensive information on substances or mixtures provided in safety data sheets. Since the provisions on safety data sheets are included in Regulation (EC) No 1907/2006 which uses the safety data sheet as the main communication tool within the supply chain of substances, it is appropriate not to duplicate the same provisions in this Regulation.

(41) To ensure proper and comprehensive information provision to consumers on the hazards and safe use of chemicals and mixtures, the use and dissemination of Internet sites and free-phone numbers should be promoted, particularly in connection with information provision on specific types of packaging.

(42) Workers and consumers worldwide would benefit from a globally harmonised hazard communication tool in the form of labelling. Therefore, the elements to be included in labels should be specified in accordance with the hazard pictograms, signal words, hazard statements and precautionary statements which form the core information of the GHS. Other information included in labels should be limited to a minimum and should not call into question the main elements.

(43) It is essential that the substances and mixtures placed on the market are well identified. However, the Agency should allow enterprises, upon their request and where necessary, to describe the chemical identity of certain substances in a way that does not put the confidential nature of their businesses at risk. Where the Agency refuses such a request, an appeal should be allowed in accordance with this Regulation. The appeal should have a suspensive effect, so that the confidential information with regard to which the request has been made, should not appear on the label while the appeal is pending.

(44) The International Union of Pure and Applied Chemistry (IUPAC) is a long-standing global authority on chemical nomenclature and terminology. Identification of substances by their IUPAC name is widespread practice worldwide and provides the standard basis for identifying substances in an international and multilingual context. It is therefore appropriate to use these names for the purposes of this Regulation.

(45) The Chemical Abstracts Service (CAS) provides a system whereby substances are added to the CAS Registry and are assigned a unique CAS Registry Number. Those CAS numbers are used in reference works, databases, and regulatory compliance documents throughout the world to identify substances without the ambiguity of chemical nomenclature. It is therefore appropriate to use the CAS numbers for the purposes of this Regulation.

(46) To limit the information on the label to the most essential information, principles of precedence should determine the most appropriate label elements for cases in which substances or mixtures possess several hazardous properties.

(47) Council Directive 91/414/EEC of 15 July 1991 concerning the placing of plant protection products on the market [1] and Directive 98/8/EC of the European Parliament and of the Council of 16 February 1998 concerning the placing of biocidal products on the market [2] should remain fully applicable to any product within their scope.

(48) Statements such as 'non-toxic', 'non-harmful', 'non-polluting', 'ecological' or other statements indicating that the substance or mixture is not hazardous or any other statements that are inconsistent with its classification should not appear on the label or packaging of any substance or mixture.

(49) In general, substances and mixtures, especially those supplied to the general public, should be supplied in packaging together with the necessary labelling information. The supply of appropriate information between professionals, including for unpackaged substances and

[1] OJ L 230, 19.8.1991, p. 1.
[2] OJ L 123, 24.4.1998, p. 1.

mixtures, is ensured by Regulation (EC) No 1907/2006. However, in exceptional circumstances, substances and mixtures may also be supplied to the general public unpackaged. Where appropriate, relevant labelling information should be supplied to the general public by other means, such as an invoice or bill.

(50) Rules for the application of labels and the location of information on labels are necessary to ensure that the information on labels can be easily understood.

(51) This Regulation should set general packaging standards, in order to ensure the safe supply of hazardous substances and mixtures.

(52) The resources of the authorities should be focused on substances of the highest concern with regard to health and to the environment. Provision should therefore be made to enable competent authorities and manufacturers, importers and downstream users to submit proposals to the Agency for a harmonised classification and labelling of substances classified for carcinogenicity, germ cell mutagenicity or reproductive toxicity categories 1A, 1B or 2, for respiratory sensitisation, or in respect of other effects on a case-by-case basis. The competent authorities of Member States should also be able to propose harmonised classification and labelling for active substances used in plant protection products and biocidal products. The Agency should give its opinion on the proposal while interested parties should have an opportunity to comment. The Commission should submit a draft decision on the final classification and labelling elements.

(53) In order to take full account of the work and experience accumulated under Directive 67/548/EEC, including the classification and labelling of specific substances listed in Annex I of Directive 67/548/EEC, all existing harmonised classifications should be converted into new harmonised classifications using the new criteria. Moreover, as the applicability of this Regulation is deferred and the harmonised classifications in accordance with the criteria of Directive 67/548/EEC are relevant for the classification of substances and mixtures during the ensuing transition period, all existing harmonised classifications should also be placed unchanged in an annex to this Regulation. By subjecting all future harmonisations of classifications to this Regulation, inconsistencies in harmonised classifications of the same substance under the existing and the new criteria should be avoided.

(54) In order to achieve the efficient functioning of the internal market for substances and mixtures, while at the same time ensuring a high level of protection for human health and the environment, rules should be established for a classification and labelling inventory. The classification and labelling for any registered or hazardous substance placed on the market should therefore be notified to the Agency to be included in the inventory.

(55) The Agency should study the possibilities for further simplification of the notification procedure in particular taking into account the needs of SMEs.

(56) Different manufacturers and importers of the same substance should make every effort to agree on a single classification for that substance except for hazard classes and differentiations subject to a harmonised classification for that substance.

(57) To ensure a harmonised level of protection for the general public, and, in particular, for persons who come into contact with certain substances, and the proper functioning of other Community legislation relying on classification and labelling, an inventory should record the classification in accordance with this Regulation agreed, if possible, by manufacturers and importers of the same substance, as well as decisions taken at Community level to harmonise the classification and labelling of some substances.

(58) The information included in the classification and labelling inventory should benefit from the same degree of accessibility and protection as that afforded by Regulation (EC) No 1907/2006, especially with regard to information which, if disclosed, risks jeopardising the commercial interests of those concerned.

(59) Member States should appoint the competent authority or competent authorities responsible for proposals for harmonised classification and labelling and the authorities responsible for the enforcement of the obligations set out in this Regulation. Member States should put in place effective monitoring and control measures in order to ensure compliance with this Regulation.

(60) It is important to provide advice to suppliers and any other interested parties, in particular SMEs, on their respective responsibilities and obligations under this Regulation. The national helpdesks already established under Regulation (EC) No 1907/2006 may act as the national helpdesks provided for under this Regulation.

(61) In order for the system established by this Regulation to operate effectively, it is important that there should be good cooperation and coordination between the Member States, the Agency and the Commission.

(62) In order to provide focal points for information on hazardous substances and mixtures, Member States should appoint bodies responsible for receiving information relating to health and to the chemical identity, components and nature of substances, including those for which the use of an alternative chemical name has been allowed in accordance with this Regulation, in addition to the competent authorities for the application and the authorities responsible for the enforcement of this Regulation.

(63) The responsible bodies, where requested by a Member State, may undertake statistical analysis to identify where improved risk management measures might be needed.

(64) Regular reports by the Member States and the Agency on the operation of this Regulation should be an indispensable means of monitoring the implementation of chemicals legislation as well as trends in this field. Conclusions drawn from findings in the reports should be useful and practical tools for reviewing the Regulation and, where necessary, for formulating proposals for amendments.

(65) The Forum for the exchange of information on enforcement in the Agency, established by Regulation (EC) No 1907/2006, should also exchange information about the enforcement of this Regulation.

(66) In order to ensure transparency, impartiality and consistency in the level of enforcement activities by Member States, it is necessary for Member States to set up an appropriate framework with a view to imposing effective, proportionate and dissuasive penalties for non-compliance with this Regulation, as non-compliance can result in damage to human health and the environment.

(67) Rules should be laid down requiring advertisements for substances meeting the criteria for classification set out in this Regulation to mention the associated hazards, in order to protect recipients of substances, including consumers. Advertisements for mixtures classified as hazardous that allow a member of the general public to conclude a contract for purchase without first having sight of the label should mention the type or types of hazard indicated on the label, for the same reason.

(68) A safeguard clause should be provided to address situations where a substance or a mixture constitutes a serious risk to human health or the environment, even if, in compliance with this Regulation, it is not classified as hazardous. Should such a situation occur, action at the UN level may be necessary in view of the global nature of trade in substances and mixtures.

(69) While many of the obligations on enterprises laid down in Regulation (EC) No 1907/2006 are triggered by classification, this Regulation should not alter the scope and impact of that Regulation, except for its provisions on safety data sheets. To ensure this, that Regulation should be amended accordingly.

(70) The application of this Regulation should be staggered to allow all parties involved, authorities, enterprises as well as stakeholders, to focus resources on preparing for new duties at the right times. Therefore, and because the classification of mixtures depends on the classification of substances, the provisions for the classification of mixtures should only be applied after the reclassification of all substances. Operators should be allowed to apply the classification criteria contained in this Regulation earlier on a voluntary basis, but in that case to avoid confusion the labelling and packaging should comply with this Regulation instead of Directives 67/548/EEC or 1999/45/EC.

(71) To avoid unnecessary burdens on enterprises, substances and mixtures which are already in the supply chain when the labelling provisions of this Regulation become applicable to them may continue to be placed on the market without relabelling for a certain period of time.

(72) Since the objectives of this Regulation, namely harmonising the classification, labelling and packaging rules, providing an obligation to classify and establishing a harmonised list of substances classified at Community level as well as a classification and labelling inventory, cannot be sufficiently achieved by the Member States and can therefore be better achieved at Community level, the Community may adopt measures, in accordance with the principle of subsidiarity as set out in Article 5 of the Treaty. In accordance with the principle of proportionality, as set out in that Article, this Regulation does not go beyond what is necessary in order to achieve those objectives.

(73) This Regulation observes the fundamental rights and principles which are acknowledged in particular in the Charter of Fundamental Rights of the European Union [1].

(74) This Regulation should contribute to the fulfilment of the Strategic Approach to International Chemical Management (SAICM) adopted on 6 February 2006 in Dubai.

(75) Subject to developments at UN level, the classification and labelling of persistent, bioaccumulative and toxic (PBT) and very persistent and very bioaccumulative (vPvB) substances should be included in this Regulation at a later stage.

(76) The measures necessary for the implementation of this Regulation should be adopted in accordance with Council Decision 1999/468/EC of 28 June 1999 laying down the procedures for the exercise of implementing powers conferred on the Commission [2].

(77) In particular, the Commission should be empowered to adapt this Regulation to technical and scientific progress, including incorporating amendments made at UN level to the GHS, in particular any such UN amendments relating to the use of information on similar mixtures. In carrying out such adaptations to technical and scientific progress the biannual working rhythm at UN level should be taken into account. Furthermore, the Commission should be empowered to decide on the harmonised classification and labelling of specific substances. Since those measures are of general scope and are designed to amend non-essential elements of this Regulation, they must be adopted in accordance with the regulatory procedure with scrutiny provided for in Article 5a of Decision 1999/468/EC.

[1] OJ C 364, 18.12.2000, p. 1.
[2] OJ L 184, 17.7.1999, p. 23.

(78) When, on imperative grounds of urgency, the normal time limits for the regulatory procedure with scrutiny cannot be complied with, the Commission should be able to apply the urgency procedure provided for in Article 5a(6) of Decision 1999/468/EC for the adoption of adaptations to technical progress.

(79) The Commission should also for the purposes of this Regulation be assisted by the Committee established by Regulation (EC) No 1907/2006, with a view to ensuring a consistent approach to the updating of chemicals legislation,

HAVE ADOPTED THIS REGULATION:

TITLE I

GENERAL ISSUES

Article 1

Purpose and scope

1. The purpose of this Regulation is to ensure a high level of protection of human health and the environment as well as the free movement of substances, mixtures and articles as referred to in Article 4(8) by:

(a) harmonising the criteria for classification of substances and mixtures, and the rules on labelling and packaging for hazardous substances and mixtures;

(b) providing an obligation for:

(i) manufacturers, importers and downstream users to classify substances and mixtures placed on the market;

(ii) suppliers to label and package substances and mixtures placed on the market;

(iii) manufacturers, producers of articles and importers to classify those substances not placed on the market that are subject to registration or notification under Regulation (EC) No 1907/2006;

(c) providing an obligation for manufacturers and importers of substances to notify the Agency of such classifications and label elements if these have not been submitted to the Agency as part of a registration under Regulation (EC) No 1907/2006;

(d) establishing a list of substances with their harmonised classifications and labelling elements at Community level in Part 3 of Annex VI;

(e) establishing a classification and labelling inventory of substances, which is made up of all notifications, submissions and harmonised classifications and labelling elements referred to in points (c) and (d).

2. This Regulation shall not apply to the following:

(a) radioactive substances and mixtures within the scope of Council Directive 96/29/Euratom of 13 May 1996 laying down basic safety standards for the protection of the health of workers and the general public against the danger arising from ionising radiation ([1]);

(b) substances and mixtures which are subject to customs supervision, provided that they do not undergo any treatment or processing, and which are in temporary storage, or in a free zone or free warehouse with a view to re-exportation, or in transit;

(c) non-isolated intermediates;

(d) substances and mixtures for scientific research and development, which are not placed on the market, provided they are used under controlled conditions in accordance with Community workplace and environmental legislation.

3. Waste as defined in Directive 2006/12/EC of the European Parliament and of the Council of 5 April 2006 on waste ([2]) is not a substance, mixture or article within the meaning of Article 2 of this Regulation.

4. Member States may allow for exemptions from this Regulation in specific cases for certain substances or mixtures, where necessary in the interests of defence.

5. This Regulation shall not apply to substances and mixtures in the following forms, which are in the finished state, intended for the final user:

(a) medicinal products as defined in Directive 2001/83/EC;

(b) veterinary medicinal products as defined in Directive 2001/82/EC;

(c) cosmetic products as defined in Directive 76/768/EEC;

(d) medical devices as defined in Directives 90/385/EEC and 93/42/EEC, which are invasive or used in direct physical contact with the human body, and in Directive 98/79/EC;

(e) food or feeding stuffs as defined in Regulation (EC) No 178/2002 including when they are used:

(i) as a food additive in foodstuffs within the scope of Directive 89/107/EEC;

(ii) as a flavouring in foodstuffs within the scope of Directive 88/388/EEC and Decision 1999/217/EC;

([1]) OJ L 159, 29.6.1996, p. 1.
([2]) OJ L 114, 27.4.2006, p. 9.

(iii) as an additive in feeding stuffs within the scope of Regulation (EC) No 1831/2003;

(iv) in animal nutrition within the scope of Directive 82/471/EEC.

6. Save where Article 33 applies this Regulation shall not apply to the transport of dangerous goods by air, sea, road, rail or inland waterways.

Article 2

Definitions

For the purpose of this Regulation, the following definitions shall apply:

1. 'hazard class' means the nature of the physical, health or environmental hazard;

2. 'hazard category' means the division of criteria within each hazard class, specifying hazard severity;

3. 'hazard pictogram' means a graphical composition that includes a symbol plus other graphic elements, such as a border, background pattern or colour that is intended to convey specific information on the hazard concerned;

4. 'signal word' means a word that indicates the relative level of severity of hazards to alert the reader to a potential hazard; the following two levels are distinguished:

   (a) 'Danger' means a signal word indicating the more severe hazard categories;

   (b) 'Warning' means a signal word indicating the less severe hazard categories;

5. 'hazard statement' means a phrase assigned to a hazard class and category that describes the nature of the hazards of a hazardous substance or mixture, including, where appropriate, the degree of hazard;

6. 'precautionary statement' means a phrase that describes recommended measure(s) to minimise or prevent adverse effects resulting from exposure to a hazardous substance or mixture due to its use or disposal;

7. 'substance' means a chemical element and its compounds in the natural state or obtained by any manufacturing process, including any additive necessary to preserve its stability and any impurity deriving from the process used, but excluding any solvent which may be separated without affecting the stability of the substance or changing its composition;

8. 'mixture' means a mixture or solution composed of two or more substances;

9. 'article' means an object which during production is given a special shape, surface or design which determines its function to a greater degree than does its chemical composition;

10. 'producer of an article' means any natural or legal person who makes or assembles an article within the Community;

11. 'polymer' means a substance consisting of molecules characterised by the sequence of one or more types of monomer units. Such molecules must be distributed over a range of molecular weights wherein differences in the molecular weight are primarily attributable to differences in the number of monomer units. A polymer comprises the following:

   (a) a simple weight majority of molecules containing at least three monomer units which are covalently bound to at least one other monomer unit or other reactant;

   (b) less than a simple weight majority of molecules of the same molecular weight.

   In the context of this definition a 'monomer unit' means the reacted form of a monomer substance in a polymer;

12. 'monomer' means a substance which is capable of forming covalent bonds with a sequence of additional like or unlike molecules under the conditions of the relevant polymer-forming reaction used for the particular process;

13. 'registrant' means the manufacturer or the importer of a substance or the producer or importer of an article submitting a registration for a substance under Regulation (EC) No 1907/2006;

14. 'manufacturing' means production or extraction of substances in the natural state;

15. 'manufacturer' means any natural or legal person established within the Community who manufactures a substance within the Community;

16. 'import' means the physical introduction into the customs territory of the Community;

17. 'importer' means any natural or legal person established within the Community who is responsible for import;

18. 'placing on the market' means supplying or making available, whether in return for payment or free of charge, to a third party. Import shall be deemed to be placing on the market;

19. 'downstream user' means any natural or legal person established within the Community, other than the manufacturer or the importer, who uses a substance, either on its own or in a mixture, in the course of his industrial or professional activities. A distributor or a consumer is not a downstream user. A re-importer exempted pursuant to Article 2(7)(c) of Regulation (EC) No 1907/2006 shall be regarded as a downstream user;

20. 'distributor' means any natural or legal person established within the Community, including a retailer, who only stores and places on the market a substance, on its own or in a mixture, for third parties;

21. 'intermediate' means a substance that is manufactured for and consumed in or used for chemical processing in order to be transformed into another substance (hereinafter referred to as 'synthesis');

22. 'non-isolated intermediate' means an intermediate that during synthesis is not intentionally removed (except for sampling) from the equipment in which the synthesis takes place. Such equipment includes the reaction vessel, its ancillary equipment, and any equipment through which the substance(s) pass(es) during a continuous flow or batch process as well as the pipework for transfer from one vessel to another for the purpose of the next reaction step, but it excludes tanks or other vessels in which the substance(s) are stored after the manufacture;

23. 'the Agency' means the European Chemicals Agency established by Regulation (EC) No 1907/2006;

24. 'competent authority' means the authority or authorities or bodies established by the Member States to carry out the obligations arising from this Regulation;

25. 'use' means any processing, formulation, consumption, storage, keeping, treatment, filling into containers, transfer from one container to another, mixing, production of an article or any other utilisation;

26. 'supplier' means any manufacturer, importer, downstream user or distributor placing on the market a substance, on its own or in a mixture, or a mixture;

27. 'alloy' means a metallic material, homogeneous on a macroscopic scale, consisting of two or more elements so combined that they cannot be readily separated by mechanical means; alloys are considered to be mixtures for the purposes of this Regulation;

28. 'UN RTDG' means the United Nations Recommendations on the Transport of Dangerous Goods;

29. 'notifier' means the manufacturer or the importer, or group of manufacturers or importers notifying to the Agency;

30. 'scientific research and development' means any scientific experimentation, analysis or chemical research carried out under controlled conditions;

31. 'cut-off value' means a threshold of any classified impurity, additive or individual constituent in a substance or in a mixture, above which threshold these shall be taken into account for determining if the substance or the mixture, respectively, shall be classified;

32. 'concentration limit' means a threshold of any classified impurity, additive or individual constituent in a substance or in a mixture that may trigger classification of the substance or the mixture, respectively;

33. 'differentiation' means distinction within hazard classes depending on the route of exposure or the nature of the effects;

34. 'M-factor' means a multiplying factor. It is applied to the concentration of a substance classified as hazardous to the aquatic environment acute category 1 or chronic category 1, and is used to derive by the summation method the classification of a mixture in which the substance is present;

35. 'package' means the complete product of the packing operation, consisting of the packaging and its contents;

36. 'packaging' means one or more receptacles and any other components or materials necessary for the receptacles to perform their containment and other safety functions;

37. 'intermediate packaging' means packaging placed between inner packaging, or articles, and outer packaging.

Article 3

Hazardous substances and mixtures and specification of hazard classes

A substance or a mixture fulfilling the criteria relating to physical hazards, health hazards or environmental hazards, laid down in Parts 2 to 5 of Annex I is hazardous and shall be classified in relation to the respective hazard classes provided for in that Annex.

Where, in Annex I, hazard classes are differentiated on the basis of the route of exposure or the nature of the effects, the substance or mixture shall be classified in accordance with such differentiation.

Article 4

General obligations to classify, label and package

1. Manufacturers, importers and downstream users shall classify substances or mixtures in accordance with Title II before placing them on the market.

2. Without prejudice to the requirements of paragraph 1, manufacturers, producers of articles and importers shall classify those substances not placed on the market in accordance with Title II where:

(a) Articles 6, 7(1) or (5), 17 or 18 of Regulation (EC) No 1907/2006 provide for registration of a substance;

(b) Articles 7(2) or 9 of Regulation (EC) No 1907/2006 provide for notification.

3. If a substance is subject to harmonised classification and labelling in accordance with Title V through an entry in Part 3 of Annex VI, that substance shall be classified in accordance with that entry, and a classification of that substance in accordance with Title II shall not be performed for the hazard classes or differentiations covered by that entry.

However, where the substance also falls within one or more hazard classes or differentiations not covered by an entry in Part 3 of Annex VI, classification under Title II shall be carried out for those hazard classes or differentiations.

4. Where a substance or mixture is classified as hazardous, suppliers shall ensure that the substance or mixture is labelled and packaged in accordance with Titles III and IV, before placing it on the market.

5. In fulfilling their responsibilities under paragraph 4, distributors may use the classification for a substance or mixture derived in accordance with Title II by an actor in the supply chain.

6. In fulfilling their responsibilities under paragraphs 1 and 4, downstream users may use the classification of a substance or mixture derived in accordance with Title II by an actor in the supply chain, provided that they do not change the composition of the substance or mixture.

7. A mixture referred to in Part 2 of Annex II that contains any substance classified as hazardous shall not be placed on the market, unless it is labelled in accordance with Title III.

8. For the purposes of this Regulation, the articles referred to in section 2.1 of Annex I shall be classified, labelled and packaged in accordance with the rules for substances and mixtures before being placed on the market.

9. Suppliers in a supply chain shall cooperate to meet the requirements for classification, labelling and packaging in this Regulation.

10. Substances and mixtures shall not be placed on the market unless they comply with this Regulation.

TITLE II

HAZARD CLASSIFICATION

CHAPTER 1

Identification and examination of information

Article 5

Identification and examination of available information on substances

1. Manufacturers, importers and downstream users of a substance shall identify the relevant available information for the purposes of determining whether the substance entails a physical, health or environmental hazard as set out in Annex I, and, in particular, the following:

(a) data generated in accordance with any of the methods referred to in Article 8(3);

(b) epidemiological data and experience on the effects on humans, such as occupational data and data from accident databases;

(c) any other information generated in accordance with section 1 of Annex XI to Regulation (EC) No 1907/2006;

(d) any new scientific information;

(e) any other information generated under internationally recognised chemical programmes.

The information shall relate to the forms or physical states in which the substance is placed on the market and in which it can reasonably be expected to be used.

2. Manufacturers, importers and downstream users shall examine the information referred to in paragraph 1 to ascertain whether it is adequate, reliable and scientifically valid for the purpose of the evaluation pursuant to Chapter 2 of this Title.

Article 6

Identification and examination of available information on mixtures

1. Manufacturers, importers and downstream users of a mixture shall identify the relevant available information on the mixture itself or the substances contained in it for the purposes of determining whether the mixture entails a physical, health or environmental hazard as set out in Annex I, and, in particular, the following:

(a) data generated in accordance with any of the methods referred to in Article 8(3) on the mixture itself or the substances contained in it;

(b) epidemiological data and experience on the effects on humans for the mixture itself or the substances contained in it, such as occupational data or data from accident databases;

(c) any other information generated in accordance with section 1 of Annex XI to Regulation (EC) No 1907/2006 for the mixture itself or the substances contained in it;

(d) any other information generated under internationally recognised chemical programmes for the mixture itself or the substances contained in it.

The information shall relate to the forms or physical states in which the mixture is placed on the market and, when relevant, in which it can reasonably be expected to be used.

2. Subject to paragraphs 3 and 4, where the information referred to in paragraph 1 is available for the mixture itself, and the manufacturer, importer or downstream user has ascertained that information to be adequate and reliable and where applicable, scientifically valid, that manufacturer, importer or downstream user shall use that information for the purposes of the evaluation pursuant to Chapter 2 of this Title.

3. For the evaluation of mixtures pursuant to Chapter 2 of this Title in relation to the 'germ cell mutagenicity', 'carcinogenicity' and 'reproductive toxicity' hazard classes referred to in sections 3.5.3.1, 3.6.3.1 and 3.7.3.1 of Annex I, the manufacturer, importer or downstream user shall only use the relevant available information referred to in paragraph 1 for the substances in the mixture.

Further, in cases where the available test data on the mixture itself demonstrate germ cell mutagenic, carcinogenic or toxic to reproduction effects which have not been identified from the information on the individual substances, those data shall also be taken into account.

4. For the evaluation of mixtures pursuant to Chapter 2 of this Title in relation to the 'biodegradation and bioaccumulation' properties within the 'hazardous to the aquatic environment' hazard class referred to in sections 4.1.2.8 and 4.1.2.9 of Annex I, the manufacturer, importer or downstream user shall only use the relevant available information referred to in paragraph 1 for the substances in the mixture.

5. Where no or inadequate test data on the mixture itself of the kind referred to in paragraph 1 are available, the manufacturer, importer or downstream user shall use other available information on individual substances and similar tested mixtures which may also be considered relevant for the purposes of determining whether the mixture is hazardous, provided that that manufacturer, importer or downstream user has ascertained that information to be adequate and reliable for the purpose of the evaluation pursuant to Article 9(4).

Article 7

Animal and human testing

1. Where new tests are carried out for the purposes of this Regulation, tests on animals within the meaning of Directive 86/609/EEC shall be undertaken only where no other alternatives, which provide adequate reliability and quality of data, are possible.

2. Tests on non-human primates shall be prohibited for the purposes of this Regulation.

3. Tests on humans shall not be performed for the purposes of this Regulation. Data obtained from other sources, such as clinical studies, can however be used for the purposes of this Regulation.

Article 8

Generating new information for substances and mixtures

1. For the purposes of determining whether a substance or a mixture entails a health or environmental hazard as set out in Annex I to this Regulation, the manufacturer, importer or downstream user may, provided that he has exhausted all other means of generating information including by applying the rules provided for in section 1 of Annex XI to Regulation (EC) No 1907/2006, perform new tests.

2. For the purposes of determining whether a substance or a mixture entails any of the physical hazards referred to in Part 2 of Annex I, the manufacturer, importer or downstream user shall perform the tests required in that Part, unless there is adequate and reliable information already available.

3. The tests referred to in paragraph 1 shall be conducted in accordance with one of the following methods:

(a) the test methods referred to in Article 13(3) of Regulation (EC) No 1907/2006;

or

(b) sound scientific principles that are internationally recognised or methods validated according to international procedures.

4. Where the manufacturer, importer or downstream user carries out new ecotoxicological or toxicological tests and analyses, these shall be carried out in compliance with Article 13(4) of Regulation (EC) No 1907/2006.

5. Where new tests for physical hazards are carried out for the purposes of this Regulation, they shall be carried out, at the latest from 1 January 2014, in compliance with a relevant recognised quality system or by laboratories complying with a relevant recognised standard.

6. Tests that are carried out for the purposes of this Regulation shall be carried out on the substance or on the mixture in the form(s) or physical state(s) in which the substance or mixture is placed on the market and in which it can reasonably be expected to be used.

CHAPTER 2

Evaluation of hazard information and decision on classification

Article 9

Evaluation of hazard information for substances and mixtures

1. Manufacturers, importers and downstream users of a substance or a mixture shall evaluate the information identified in accordance with Chapter 1 of this Title by applying to it the criteria for classification for each hazard class or differentiation in Parts 2 to 5 of Annex I, so as to ascertain the hazards associated with the substance or mixture.

2. In evaluating available test data for a substance or a mixture which have been obtained from test methods other than those referred to in Article 8(3), manufacturers, importers and downstream users shall compare the test methods employed with those indicated in that Article in order to determine whether the use of those test methods affects the evaluation referred to in paragraph 1 of this Article.

3. Where the criteria cannot be applied directly to available identified information, manufacturers, importers and downstream users shall carry out an evaluation by applying a weight of evidence determination using expert judgement in accordance with section 1.1.1 of Annex I to this Regulation, weighing all available information having a bearing on the determination of the hazards of the substance or the mixture, and in accordance with section 1.2 of Annex XI to Regulation (EC) No 1907/2006.

4. Where only the information referred to in Article 6(5) is available, manufacturers, importers and downstream users shall apply the bridging principles referred to in section 1.1.3 and in each section of Parts 3 and 4 of Annex I for the purposes of the evaluation.

However, where that information permits the application neither of the bridging principles nor the principles for using expert judgement and weight of evidence determination as described in Part 1 of Annex I, manufacturers, importers and downstream users shall evaluate the information by applying the other method or methods described in each section of Parts 3 and 4 of Annex I.

5. When evaluating the available information for the purposes of classification, the manufacturers, importers and downstream users shall consider the forms or physical states in which the substance or mixture is placed on the market and in which it can reasonably be expected to be used.

Article 10

Concentration limits and M-factors for classification of substances and mixtures

1. Specific concentration limits and generic concentration limits are limits assigned to a substance indicating a threshold at or above which the presence of that substance in another substance or in a mixture as an identified impurity, additive or individual constituent leads to the classification of the substance or mixture as hazardous.

Specific concentration limits shall be set by the manufacturer, importer or downstream user where adequate and reliable scientific information shows that the hazard of a substance is evident when the substance is present at a level below the concentrations set for any hazard class in Part 2 of Annex I or below the generic concentration limits set for any hazard class in Parts 3, 4 and 5 of Annex I.

In exceptional circumstances specific concentration limits may be set by the manufacturer, importer or downstream user where he has adequate, reliable and conclusive scientific information that a hazard of a substance classified as hazardous is not evident at a level above the concentrations set for the relevant hazard class in Part 2 of Annex I or above the generic concentration limits set for the relevant hazard class in Parts 3, 4 and 5 of that Annex.

2. M-factors for substances classified as hazardous to the aquatic environment, acute category 1 or chronic category 1, shall be established by manufacturers, importers and downstream users.

3. Notwithstanding paragraph 1, specific concentration limits shall not be set for harmonised hazard classes or differentiations for substances included in Part 3 of Annex VI.

4. Notwithstanding paragraph 2, M-factors shall not be set for harmonised hazard classes or differentiations for substances included in Part 3 of Annex VI for which an M-factor is given in that Part.

However, where an M-factor is not given in Part 3 of Annex VI for substances classified as hazardous to the aquatic environment, acute category 1 or chronic category 1, an M-factor based on available data for the substance shall be set by the manufacturer, importer or downstream user. When a mixture including the substance is classified by the manufacturer, importer or downstream user using the summation method, this M-factor shall be used.

5. In setting the specific concentration limit or M-factor manufacturers, importers and downstream users shall take into account any specific concentration limits or M-factors for that substance which have been included in the classification and labelling inventory.

6. Specific concentration limits set in accordance with paragraph 1 shall take precedence over the concentrations in the relevant sections of Part 2 of Annex I or the generic concentration limits for classification in the relevant sections of Parts 3, 4 and 5 of Annex I.

7. The Agency shall provide further guidance for the application of paragraphs 1 and 2.

Article 11

Cut-off values

1. Where a substance contains another substance, itself classified as hazardous, whether in the form of an identified impurity, additive or individual constituent, this shall be taken into account for the purposes of classification, if the concentration of the identified impurity, additive or individual constituent is equal to, or greater than, the applicable cut-off value in accordance with paragraph 3.

2. Where a mixture contains a substance classified as hazardous, whether as a component or in the form of an identified impurity or additive, this information shall be taken into account for the purposes of classification, if the concentration of that substance is equal to or greater than its cut-off value in accordance with paragraph 3.

3. The cut-off value referred to in paragraphs 1 and 2 shall be determined as set out in section 1.1.2.2 of Annex I.

Article 12

Specific cases requiring further evaluation

Where, as a result of the evaluation carried out pursuant to Article 9, the following properties or effects are identified, manufacturers, importers and downstream users shall take them into account for the purposes of classification:

(a) adequate and reliable information demonstrates that in practice the physical hazards of a substance or a mixture differ from those shown by tests;

(b) conclusive scientific experimental data show that the substance or mixture is not biologically available and those data have been ascertained to be adequate and reliable;

(c) adequate and reliable scientific information demonstrates the potential occurrence of synergistic or antagonistic effects among the substances in a mixture for which the evaluation was decided on the basis of the information for the substances in the mixture.

Article 13

Decision to classify substances and mixtures

If the evaluation undertaken pursuant to Article 9 and Article 12 shows that the hazards associated with the substance or mixture meet the criteria for classification in one or more hazard classes or differentiations in Parts 2 to 5 of Annex I, manufacturers, importers and downstream users shall classify the substance or mixture in relation to the relevant hazard class or classes or differentiations by assigning the following:

(a) one or more hazard categories for each relevant hazard class or differentiation;

(b) subject to Article 21, one or more hazard statements corresponding to each hazard category assigned in accordance with (a).

Article 14

Specific rules for the classification of mixtures

1. The classification of a mixture shall not be affected where the evaluation of the information indicates any of the following:

(a) that the substances in the mixture react slowly with atmospheric gases, in particular oxygen, carbon dioxide, water vapour, to form different substances at low concentration;

(b) that the substances in the mixture react very slowly with other substances in the mixture to form different substances at low concentration;

(c) that the substances in the mixture may self-polymerise to form oligomers or polymers, at low concentration.

2. A mixture need not be classified for explosive, oxidising, or flammable properties as referred to in Part 2 of Annex I provided that any of the following requirements are met:

(a) none of the substances in the mixture possesses any of those properties and, on the basis of the information available to the supplier, the mixture is unlikely to present hazards of this kind;

(b) in the event of a change in the composition of a mixture, scientific evidence indicates that an evaluation of the information on the mixture will not lead to a change in classification;

(c) where a mixture is placed on the market in the form of an aerosol dispenser, it satisfies Article 8(1a) of Council Directive 75/324/EEC of 20 May 1975 on the approximation of the laws of the Member States relating to aerosol dispensers [1].

[1] OJ L 147, 9.6.1975, p. 40.

Article 15

Review of classification for substances and mixtures

1. Manufacturers, importers and downstream users shall take all reasonable steps available to them to make themselves aware of new scientific or technical information that may affect the classification of the substances or mixtures they place on the market. When a manufacturer, importer or downstream user becomes aware of such information which he considers to be adequate and reliable, that manufacturer, importer or downstream user shall without undue delay carry out a new evaluation in accordance with this Chapter.

2. Where the manufacturer, importer or downstream user introduces a change to a mixture that has been classified as hazardous, that manufacturer, importer or downstream user shall carry out a new evaluation in accordance with this Chapter where the change is either of the following:

(a) a change in the composition of the initial concentration of one or more of the hazardous constituents in concentrations at or above the limits in Table 1.2 of Part 1 of Annex I;

(b) a change in the composition involving the substitution or addition of one or more constituents in concentrations at or above the cut-off value referred to in Article 11(3).

3. A new evaluation in accordance with paragraphs 1 and 2 shall not be required if there is valid scientific justification that this will not result in a change of classification.

4. Manufacturers, importers and downstream users shall adapt the classification of the substance or the mixture in accordance with the results of the new evaluation except where there are harmonised hazard classes or differentiations for substances included in Part 3 of Annex VI.

5. For paragraphs 1 to 4 of this Article, when the substance or mixture concerned is within the scope of Directive 91/414/EEC or Directive 98/8/EC, the requirements of those Directives shall also apply.

Article 16

Classification of substances included in the classification and labelling inventory

1. Manufacturers and importers may classify a substance differently from the classification already included in the classification and labelling inventory, provided they submit the reasons for the classification to the Agency together with the notification in accordance with Article 40.

2. Paragraph 1 shall not apply if the classification included in the classification and labelling inventory is a harmonised classification included in Part 3 of Annex VI.

TITLE III

HAZARD COMMUNICATION IN THE FORM OF LABELLING

CHAPTER 1

Content of the label

Article 17

General rules

1. A substance or mixture classified as hazardous and contained in packaging shall bear a label including the following elements:

(a) the name, address and telephone number of the supplier(s);

(b) the nominal quantity of the substance or mixture in the package made available to the general public, unless this quantity is specified elsewhere on the package;

(c) product identifiers as specified in Article 18;

(d) where applicable, hazard pictograms in accordance with Article 19;

(e) where applicable, signal words in accordance with Article 20;

(f) where applicable, hazard statements in accordance with Article 21;

(g) where applicable, the appropriate precautionary statements in accordance with Article 22;

(h) where applicable, a section for supplemental information in accordance with Article 25.

2. The label shall be written in the official language(s) of the Member State(s) where the substance or mixture is placed on the market, unless the Member State(s) concerned provide(s) otherwise.

Suppliers may use more languages on their labels than those required by the Member States, provided that the same details appear in all languages used.

Article 18

Product identifiers

1. The label shall include details permitting the identification of the substance or mixture (hereinafter referred to as 'product identifiers').

The term used for identification of the substance or mixture shall be the same as that used in the safety data sheet drawn up in accordance with Article 31 of Regulation (EC) No 1907/2006 (hereinafter referred to as 'safety data sheet'), without prejudice to Article 17(2) of this Regulation.

2. The product identifier for a substance shall consist of at least the following:

(a) if the substance is included in Part 3 of Annex VI, a name and an identification number as given therein;

(b) if the substance is not included in Part 3 of Annex VI, but appears in the classification and labelling inventory, a name and an identification number as given therein;

(c) if the substance is not included in Part 3 of Annex VI nor in the classification and labelling inventory, the number provided by the CAS (hereinafter referred to as 'the CAS number'), together with the name set out in the nomenclature provided by the IUPAC (hereinafter referred to as 'the IUPAC Nomenclature'), or the CAS number together with another international chemical name(s); or

(d) if the CAS number is not available, the name set out in the IUPAC Nomenclature or another international chemical name(s).

Where the name in the IUPAC nomenclature exceeds 100 characters, one of the other names (usual name, trade name, abbreviation) referred to in section 2.1.2 of Annex VI to Regulation (EC) No 1907/2006 may be used provided that the notification in accordance with Article 40 includes both the name set out in the IUPAC Nomenclature and the other name used.

3. The product identifier for a mixture shall consist of both of the following:

(a) the trade name or the designation of the mixture;

(b) the identity of all substances in the mixture that contribute to the classification of the mixture as regards acute toxicity, skin corrosion or serious eye damage, germ cell mutagenicity, carcinogenicity, reproductive toxicity, respiratory or skin sensitisation, specific target organ toxicity (STOT) or aspiration hazard.

Where, in the case referred to in (b), that requirement leads to the provision of multiple chemical names, a maximum of four chemical names shall suffice, unless more than four names are needed to reflect the nature and the severity of the hazards.

The chemical names selected shall identify the substances primarily responsible for the major health hazards which have given rise to the classification and the choice of the corresponding hazard statements.

Article 19

Hazard pictograms

1. The label shall include the relevant hazard pictogram(s), intended to convey specific information on the hazard concerned.

2. Subject to Article 33, hazard pictograms shall fulfil the requirements laid down in section 1.2.1 of Annex I and in Annex V.

3. The hazard pictogram relevant for each specific classification is set out in the tables indicating the label elements required for each hazard class in Annex I.

Article 20

Signal words

1. The label shall include the relevant signal word in accordance with the classification of the hazardous substance or mixture.

2. The signal word relevant for each specific classification is set out in the tables indicating the label elements required for each hazard class in Parts 2 to 5 of Annex I.

3. Where the signal word 'Danger' is used on the label, the signal word 'Warning' shall not appear on the label.

Article 21

Hazard statements

1. The label shall include the relevant hazard statements in accordance with the classification of the hazardous substance or mixture.

2. The hazard statements relevant for each classification are set out in the tables indicating the label elements required for each hazard class in Parts 2 to 5 of Annex I.

3. Where a substance is included in Part 3 of Annex VI, the hazard statement relevant for each specific classification covered by the entry in that Part shall be used on the label, together with the hazard statements referred to in paragraph 2 for any other classification not covered by that entry.

4. The hazard statements shall be worded in accordance with Annex III.

Article 22

Precautionary statements

1. The label shall include the relevant precautionary statements.

2. The precautionary statements shall be selected from those set out in the tables in Parts 2 to 5 of Annex I indicating the label elements for each hazard class.

3. The precautionary statements shall be selected in accordance with the criteria laid down in Part 1 of Annex IV taking into account the hazard statements and the intended or identified use or uses of the substance or the mixture.

4. The precautionary statements shall be worded in accordance with Part 2 of Annex IV.

Article 23

Derogations from labelling requirements for special cases

The specific provisions on labelling laid down in section 1.3 of Annex I shall apply in respect of the following:

(a) transportable gas cylinders;

(b) gas containers intended for propane, butane or liquefied petroleum gas;

(c) aerosols and containers fitted with a sealed spray attachment and containing substances or mixtures classified as presenting an aspiration hazard;

(d) metals in massive form, alloys, mixtures containing polymers, mixtures containing elastomers;

(e) explosives, as referred to in section 2.1 of Annex I, placed on the market with a view to obtaining an explosive or pyrotechnic effect.

Article 24

Request for use of an alternative chemical name

1. The manufacturer, importer or downstream user of a substance in a mixture may submit a request to the Agency to use an alternative chemical name which refers to that substance in a mixture either by means of a name that identifies the most important functional chemical groups or by means of an alternative designation, where the substance meets the criteria set out in Part 1 of Annex I and where he can demonstrate that disclosure on the label or in the safety data sheet of the chemical identity of that substance puts the confidential nature of his business, in particular his intellectual property rights, at risk.

2. Any request referred to in paragraph 1 of this Article shall be made in the format referred to in Article 111 of Regulation (EC) No 1907/2006 and shall be accompanied by a fee.

The level of the fees shall be determined by the Commission in accordance with the regulatory procedure referred to in Article 54(2) of this Regulation.

A reduced fee shall be set for SMEs.

3. The Agency may require further information from the manufacturer, importer or downstream user making the request if such information is necessary to take a decision. If the Agency raises no objection within six weeks of the request or the receipt of further required information, the use of the requested name shall be deemed to be allowed.

4. If the Agency does not accept the request, the practical arrangements referred to in Article 118(3) of Regulation (EC) No 1907/2006 shall apply.

5. The Agency shall inform competent authorities of the outcome of the request in accordance with paragraph 3 or 4 and provide them with the information submitted by the manufacturer, importer or downstream user.

6. Where new information shows that an alternative chemical name used does not provide sufficient information for necessary health and safety precautions to be taken at the workplace and to ensure that risks from handling the mixture can be controlled, the Agency shall review its decision on the use of that alternative chemical name. The Agency may withdraw its decision or amend it by a decision specifying which alternative chemical name is allowed to be used. If the Agency withdraws or amends its decision, the practical arrangements referred to in Article 118(3) of Regulation (EC) No 1907/2006 shall apply.

7. Where the use of an alternative chemical name has been allowed, but the classification of the substance in a mixture for which the alternative name is used no longer meets the criteria set out in section 1.4.1 of Annex I, the supplier of that substance in a mixture shall use the product identifier for the substance in accordance with Article 18 on the label and in the safety data sheet, and not the alternative chemical name.

8. For substances, whether on their own or in a mixture, where a justification in accordance with Article 10(a)(xi) of Regulation (EC) No 1907/2006 regarding information referred to in Article 119(2)(f) or (g) of that Regulation has been accepted as valid by the Agency, the manufacturer, importer or downstream user may use on the label and in the safety data sheet a name that will be made publicly available over the Internet. For those substances in a mixture for which Article 119(2)(f) or (g) of that Regulation no longer applies, the manufacturer, importer or downstream user may submit a request to the Agency to use an alternative chemical name as provided for in paragraph 1 of this Article.

9. Where the supplier of a mixture, before 1 June 2015, has demonstrated under Article 15 of Directive 1999/45/EC that the disclosure of the chemical identity of a substance in a mixture puts the confidential nature of his business at risk, he can continue to use the agreed alternative name for the purposes of this Regulation.

Article 25

Supplemental information on the label

1. Statements shall be included in the section for supplemental information on the label where a substance or mixture classified as hazardous has the physical properties or health properties referred to in sections 1.1 and 1.2 of Annex II.

The statements shall be worded in accordance with sections 1.1 and 1.2 of Annex II and Part 2 of Annex III.

Where a substance is included in Part 3 of Annex VI, any supplemental hazard statements given therein for the substance shall be included in the supplemental information on the label.

2. A statement shall be included in the section for supplemental information on the label where a substance or mixture classified as hazardous falls within the scope of Directive 91/414/EEC.

The statement shall be worded in accordance with Part 4 of Annex II and Part 3 of Annex III to this Regulation.

3. The supplier may include supplemental information in the section for supplemental information on the label other than that referred to in paragraphs 1 and 2, provided that that information does not make it more difficult to identify the label elements referred to in Article 17(1) (a) to (g) and that it provides further details and does not contradict or cast doubt on the validity of the information specified by those elements.

4. Statements such as 'non-toxic', 'non-harmful', 'non-polluting', 'ecological' or any other statements indicating that the substance or mixture is not hazardous or any other statements that are inconsistent with the classification of that substance or mixture shall not appear on the label or packaging of any substance or mixture.

5. Where a substance or mixture is classified in accordance with Part 5 of Annex I,

(a) the hazard pictogram shall not be included on the label;

(b) the signal words, hazard statements and precautionary statements shall be placed in the supplemental information section of the label.

6. Where a mixture contains any substance classified as hazardous, it shall be labelled in accordance with Part 2 of Annex II.

The statements shall be worded in accordance with Part 3 of Annex III and shall be placed in the supplemental information section of the label.

The label shall also include the product identifier referred to in Article 18 and the name, address and telephone number of the supplier of the mixture.

Article 26

Principles of precedence for hazard pictograms

1. Where the classification of a substance or mixture would result in more than one hazard pictogram on the label, the following rules of precedence shall apply to reduce the number of hazard pictograms required:

(a) if the hazard pictogram 'GHS01' applies, the use of the hazard pictograms 'GHS02' and 'GHS03' shall be optional, except in cases where more than one of these hazard pictograms are compulsory;

(b) if the hazard pictogram 'GHS06' applies, the hazard pictogram 'GHS07' shall not appear;

(c) if the hazard pictogram 'GHS05' applies, the hazard pictogram 'GHS07' shall not appear for skin or eye irritation;

(d) if the hazard pictogram 'GHS08' applies for respiratory sensitisation, the hazard pictogram 'GHS07' shall not appear for skin sensitisation or for skin and eye irritation.

2. Where the classification of a substance or mixture would result in more than one hazard pictogram for the same hazard class the label shall include the hazard pictogram corresponding to the most severe hazard category for each hazard class concerned.

For substances that are included in Part 3 of Annex VI and also subject to classification pursuant to Title II, the label shall include the hazard pictogram corresponding to the most severe hazard category for each relevant hazard class.

Article 27

Principles of precedence for hazard statements

If a substance or mixture is classified within several hazard classes or differentiations of a hazard class, all hazard statements resulting from the classification shall appear on the label, unless there is evident duplication or redundancy.

Article 28

Principles of precedence for precautionary statements

1. Where the selection of the precautionary statements results in certain precautionary statements being clearly redundant or unnecessary given the specific substance, mixture or packaging, such statements shall be omitted from the label.

2. Where the substance or mixture is supplied to the general public, one precautionary statement addressing the disposal of that substance or mixture as well as the disposal of packaging shall appear on the label, unless not required under Article 22.

In all other cases, a precautionary statement addressing disposal shall not be required, where it is clear that the disposal of the substance or mixture or the packaging does not present a hazard to human health or the environment.

3. Not more than six precautionary statements shall appear on the label, unless necessary to reflect the nature and the severity of the hazards.

Article 29

Exemptions from labelling and packaging requirements

1. Where the packaging of a substance or a mixture is either in such a shape or form or is so small that it is impossible to meet the requirements of Article 31 for a label in the languages of the

Member State in which the substance or mixture is placed on the market, the label elements in accordance with the first subparagraph of Article 17(2) shall be provided in accordance with section 1.5.1 of Annex I.

2. If the full label information cannot be provided in the way specified in paragraph 1 the label information may be reduced in accordance with section 1.5.2 of Annex I.

3. When a hazardous substance or mixture referred to in Part 5 of Annex II is supplied to the general public without packaging it shall be accompanied by a copy of the label elements in accordance with Article 17.

4. For certain mixtures classified as hazardous to the environment, exemptions to certain provisions on environmental labelling or specific provisions in relation to environmental labelling may be determined in accordance with the procedure referred to in Article 53, where it can be demonstrated that there would be a reduction in the environmental impact. Such exemptions or specific provisions are defined in Part 2 of Annex II.

5. The Commission may request the Agency to prepare and submit to it further draft exemptions from labelling and packaging requirements.

Article 30

Updating information on labels

1. The supplier shall ensure that the label is updated, without undue delay, following any change to the classification and labelling of that substance or mixture, where the new hazard is more severe or where new supplemental labelling elements are required under Article 25, taking into account the nature of the change as regards the protection of human health and the environment. Suppliers shall cooperate in accordance with Article 4(9) to complete the changes to the labelling without undue delay.

2. Where labelling changes are required other than those referred to in paragraph 1, the supplier shall ensure that the label is updated within 18 months.

3. The supplier of a substance or a mixture within the scope of Directives 91/414/EEC or 98/8/EC shall update the label in accordance with those Directives.

CHAPTER 2

Application of labels

Article 31

General rules for the application of labels

1. Labels shall be firmly affixed to one or more surfaces of the packaging immediately containing the substance or mixture and shall be readable horizontally when the package is set down normally.

2. The colour and presentation of any label shall be such that the hazard pictogram stands out clearly.

3. The label elements referred to in Article 17(1) shall be clearly and indelibly marked. They shall stand out clearly from the background and be of such size and spacing as to be easily read.

4. The shape, colour and the size of a hazard pictogram as well as the dimensions of the label shall be as set out in section 1.2.1 of Annex I.

5. A label shall not be required when the label elements referred to in Article 17(1) are shown clearly on the packaging itself. In such cases, the requirements of this Chapter applicable to a label shall be applied to the information shown on the packaging.

Article 32

Location of information on the label

1. The hazard pictograms, signal word, hazard statements and precautionary statements shall be located together on the label.

2. The supplier may decide the order of the hazard statements on the label. However, subject to paragraph 4, all hazard statements shall be grouped on the label by language.

The supplier may decide the order of the precautionary statements on the label. However, subject to paragraph 4, all precautionary statements shall be grouped on the label by language.

3. Groups of hazard statements and groups of precautionary statements referred to in paragraph 2 shall be located together on the label by language.

4. The supplemental information shall be placed in the supplemental information section referred to in Article 25, and shall be located with the other label elements specified in Article 17(1)(a) to (g).

5. In addition to its use in hazard pictograms, colour may be used on other areas of the label to implement special labelling requirements.

6. Label elements resulting from the requirements provided for in other Community acts shall be placed in the section for supplemental information on the label referred to in Article 25.

Article 33

Specific rules for labelling of outer packaging, inner packaging and single packaging

1. Where a package consists of an outer and an inner packaging, together with any intermediate packaging, and the outer packaging meets labelling provisions in accordance with the rules on the transport of dangerous goods, the inner and any intermediate packaging shall be labelled in accordance with this Regulation. The outer packaging may also be labelled in

accordance with this Regulation. Where the hazard pictogram(s) required by this Regulation relate to the same hazard as in the rules for the transport of dangerous goods, the hazard pictogram(s) required by this Regulation need not appear on the outer packaging.

2. Where the outer packaging of a package is not required to meet labelling provisions in accordance with rules on the transport of dangerous goods, both the outer and any inner packaging, including any intermediate packaging, shall be labelled in accordance with this Regulation. However, if the outer packaging permits the inner or intermediate packaging labelling to be clearly seen, the outer packaging need not be labelled.

3. Single packages that meet the labelling provisions in accordance with the rules on the transport of dangerous goods shall be labelled both in accordance with this Regulation and the rules on the transport of dangerous goods. Where the hazard pictogram(s) required by this Regulation relate to the same hazard as in rules on the transport of dangerous goods, the hazard pictogram(s) required by this Regulation need not appear.

Article 34

Report on communication on safe use of chemicals

1. By 20 January 2012, the Agency shall carry out a study on the communication of information to the general public on the safe use of substances and mixtures and the potential need for additional information on labels. This study shall be carried out in consultation with competent authorities and stakeholders and drawing as appropriate on relevant best practice.

2. Without prejudice to the labelling rules provided for in this Title, the Commission shall, on the basis of the study referred to in paragraph 1, submit a report to the European Parliament and the Council and, if justified, present a legislative proposal to amend this Regulation.

TITLE IV

PACKAGING

Article 35

Packaging

1. Packaging containing hazardous substances or mixtures shall satisfy the following requirements:

(a) the packaging shall be designed and constructed so that its contents cannot escape, except in cases where other more specific safety devices are prescribed;

(b) the materials constituting the packaging and fastenings shall not be susceptible to damage by the contents, or liable to form hazardous compounds with the contents;

(c) the packaging and fastenings shall be strong and solid throughout to ensure that they will not loosen and will safely meet the normal stresses and strains of handling;

(d) packaging fitted with replaceable fastening devices shall be designed so that it can be refastened repeatedly without the contents escaping.

2. Packaging containing a hazardous substance or a mixture supplied to the general public shall not have either a shape or design likely to attract or arouse the active curiosity of children or to mislead consumers, or have a similar presentation or a design used for foodstuff or animal feeding stuff or medicinal or cosmetic products, which would mislead consumers.

Where the packaging contains a substance or mixture which meets the requirements in section 3.1.1 of Annex II it shall have a child-resistant fastening in accordance with sections 3.1.2, 3.1.3 and 3.1.4.2 of Annex II.

Where the packaging contains a substance or mixture which meets the requirements in section 3.2.1 of Annex II it shall bear a tactile warning of danger in accordance with section 3.2.2 of Annex II.

3. The packaging of substances and mixtures shall be deemed to satisfy the requirements of paragraph 1(a), (b) and (c) if it complies with the requirements of the rules on the transport of dangerous goods by air, sea, road, rail or inland waterways.

TITLE V

HARMONISATION OF CLASSIFICATION AND LABELLING OF SUBSTANCES AND THE CLASSIFICATION AND LABELLING INVENTORY

CHAPTER 1

Establishing harmonised classification and labelling of substances

Article 36

Harmonisation of classification and labelling of substances

1. A substance that fulfils the criteria set out in Annex I for the following shall normally be subject to harmonised classification and labelling in accordance with Article 37:

(a) respiratory sensitisation, category 1 (Annex I, section 3.4);

(b) germ cell mutagenicity, category 1A, 1B or 2 (Annex I, section 3.5);

(c) carcinogenicity, category 1A, 1B or 2 (Annex I, section 3.6);

(d) reproductive toxicity, category 1A, 1B or 2 (Annex I, section 3.7).

2. A substance that is an active substance in the meaning of Directive 91/414/EEC or Directive 98/8/EC shall normally be subject to harmonised classification and labelling. For such substances, the procedures set out in Article 37, paragraphs 1, 4, 5 and 6 shall apply.

3. Where a substance fulfils the criteria for other hazard classes or differentiations than those referred to in paragraph 1 and does not fall under paragraph 2, a harmonised classification and labelling in accordance with Article 37 may also be added to Annex VI on a case-by-case basis, if justification is provided demonstrating the need for such action at Community level.

Article 37

Procedure for harmonisation of classification and labelling of substances

1. A competent authority may submit to the Agency a proposal for harmonised classification and labelling of substances and, where appropriate, specific concentration limits or M-factors, or a proposal for a revision thereof.

The proposal shall follow the format set out in Part 2 of Annex VI and contain the relevant information provided for in Part 1 of Annex VI.

2. A manufacturer, importer or downstream user of a substance may submit to the Agency a proposal for harmonised classification and labelling of that substance and, where appropriate, specific concentration limits or M-factors, provided that there is no entry in Part 3 of Annex VI for such a substance in relation to the hazard class or differentiation covered by that proposal.

The proposal shall be drawn up in accordance with the relevant Parts of sections 1, 2 and 3 of Annex I to Regulation (EC) No 1907/2006 and it shall follow the format set out in Part B of the Chemical Safety Report of section 7 of that Annex. It shall contain the relevant information provided for in Part 1 of Annex VI to this Regulation. Article 111 of Regulation (EC) No 1907/2006 shall apply.

3. Where the proposal of the manufacturer, importer or downstream user concerns the harmonised classification and labelling of a substance in accordance with Article 36(3), it shall be accompanied by the fee determined by the Commission in accordance with the regulatory procedure referred to in Article 54(2).

4. The Committee for Risk Assessment of the Agency set up pursuant to Article 76(1)(c) of Regulation (EC) No 1907/2006 shall adopt an opinion on any proposal submitted pursuant to paragraphs 1 or 2 within 18 months of receipt of the proposal, giving the parties concerned the opportunity to comment. The Agency shall forward this opinion and any comments to the Commission.

5. Where the Commission finds that the harmonisation of the classification and labelling of the substance concerned is appropriate, it shall, without undue delay, submit a draft decision concerning the inclusion of that substance together with the relevant classification and labelling elements in Table 3.1 of Part 3 of Annex VI and, where appropriate, the specific concentration limits or M-factors.

A corresponding entry shall be included in Table 3.2 of Part 3 of Annex VI subject to the same conditions, until 31 May 2015.

That measure, designed to amend non-essential elements of this Regulation, shall be adopted in accordance with the regulatory procedure with scrutiny referred to in Article 54(3). On imperative grounds of urgency, the Commission may have recourse to the urgency procedure referred to in Article 54(4).

6. Manufacturers, importers and downstream users who have new information which may lead to a change of the harmonised classification and labelling elements of a substance in Part 3 of Annex VI shall submit a proposal in accordance with the second subparagraph of paragraph 2 to the competent authority in one of the Member States in which the substance is placed on the market.

Article 38

Content of opinions and decisions for harmonised classification and labelling in Part 3 of Annex VI; accessibility of information

1. Any opinion referred to in Article 37(4) and any decision according to Article 37(5) shall at least specify for each substance:

(a) the identity of the substance as specified in sections 2.1 to 2.3.4 of Annex VI to Regulation (EC) No 1907/2006;

(b) the classification of the substance referred to in Article 36, including a statement of reasons;

(c) the specific concentration limits or M-factors, where applicable;

(d) the label elements specified in points (d), (e) and (f) of Article 17(1) for the substance, together with any supplemental hazard statements for the substance, determined in accordance with Article 25(1);

(e) any other parameter enabling an assessment to be made of the health or environmental hazard of mixtures containing the hazardous substance in question or of substances containing such hazardous substances as identified impurities, additives and constituents, if relevant.

2. When making publicly available an opinion or a decision as referred to in Article 37(4) and (5) of this Regulation, Article 118(2) and Article 119 of Regulation (EC) No 1907/2006 shall apply.

CHAPTER 2

Classification and labelling inventory

Article 39

Scope

This Chapter shall apply to:

(a) substances subject to registration in accordance with Regulation (EC) No 1907/2006;

(b) substances within the scope of Article 1 which meet the criteria for classification as hazardous and are placed on the market either on their own or in a mixture above the concentration limits specified in this Regulation or Directive 1999/45/EC, where relevant, which results in the classification of the mixture as hazardous.

Article 40

Obligation to notify the Agency

1. Any manufacturer or importer, or group of manufacturers or importers (hereinafter referred to as 'the notifier(s)'), who places on the market a substance referred to in Article 39, shall notify to the Agency the following information in order for it to be included in the inventory referred to in Article 42:

(a) the identity of the notifier(s) responsible for placing the substance or substances on the market as specified in section 1 of Annex VI to Regulation (EC) No 1907/2006;

(b) the identity of the substance or substances as specified in section 2.1 to 2.3.4 to Annex VI to Regulation (EC) No 1907/2006;

(c) the classification of the substance or substances in accordance with Article 13;

(d) where a substance has been classified in some but not all hazard classes or differentiations, an indication of whether this is due to lack of data, inconclusive data, or data which are conclusive although insufficient for classification;

(e) specific concentration limits or M-factors, where applicable, in accordance with Article 10 of this Regulation together with a justification using the relevant Parts of sections 1, 2 and 3 of Annex I to Regulation (EC) No 1907/2006;

(f) the label elements specified in points (d), (e) and (f) of Article 17(1) for the substance or substances together with any supplemental hazard statements for the substance, determined in accordance with Article 25(1).

The information referred to in (a) to (f) shall not be notified, if it has been submitted to the Agency as part of a registration pursuant to Regulation (EC) No 1907/2006, or if it has already been notified by that notifier.

The notifier shall submit this information in the format specified pursuant to Article 111 of Regulation (EC) No 1907/2006.

2. The information listed in paragraph 1 shall be updated and notified to the Agency by the notifier(s) concerned when, pursuant to the review in Article 15(1), a decision to change the classification and labelling of the substance has been taken.

3. Substances placed on the market on or after 1 December 2010 shall be notified in accordance with paragraph 1 within one month after their placing on the market.

However, substances placed on the market before 1 December 2010 may be notified in accordance with paragraph 1 before that date.

Article 41

Agreed entries

Where the notification in Article 40(1) results in different entries on the inventory referred to in Article 42 for the same substance, the notifiers and registrants shall make every effort to come to an agreed entry to be included in the inventory. The notifiers shall inform the Agency accordingly.

Article 42

The classification and labelling inventory

1. The Agency shall establish and maintain a classification and labelling inventory in the form of a database.

The information notified pursuant to Article 40(1) shall be included in the inventory, as well as information submitted as part of registrations under Regulation (EC) No 1907/2006.

Information in the inventory which corresponds to the information referred to in Article 119(1) of Regulation (EC) No 1907/2006 shall be publicly accessible. The Agency shall grant access to the other information on each substance in the inventory to the notifiers and registrants who have submitted information on that substance in accordance with Article 29(1) of Regulation (EC) No 1907/2006. It shall grant access to such information to other parties subject to Article 118 of that Regulation.

2. The Agency shall update the inventory when it receives updated information in accordance with Article 40(2) or Article 41.

3. In addition to the information referred to in paragraph 1, the Agency shall, where applicable, include the following information in each entry:

(a) whether, in respect of the entry, there is harmonised classification and labelling at Community level by inclusion in Part 3 of Annex VI;

(b) whether, in respect of the entry, it is a joint entry between registrants of the same substance as referred to in Article 11(1) of Regulation (EC) No 1907/2006;

(c) whether it is an agreed entry of two or more notifiers or registrants in accordance with Article 41;

(d) whether the entry differs from another entry on the inventory for the same substance.

The information referred to in (a) shall be updated where a decision is taken in accordance with Article 37(5).

TITLE VI

COMPETENT AUTHORITIES AND ENFORCEMENT

Article 43

Appointment of competent authorities and enforcement authorities and cooperation between authorities

Member States shall appoint the competent authority or competent authorities responsible for proposals for harmonised classification and labelling and the authorities responsible for the enforcement of the obligations set out in this Regulation.

The competent authorities and the authorities responsible for enforcement shall cooperate with each other in the performance of their tasks under this Regulation and shall give the corresponding authorities of other Member States all necessary and useful support to this end.

Article 44

Helpdesk

Member States shall establish national helpdesks to provide advice to manufacturers, importers, distributors, downstream users and any other interested parties on their respective responsibilities and obligations under this Regulation.

Article 45

Appointment of bodies responsible for receiving information relating to emergency health response

1. Member States shall appoint a body or bodies responsible for receiving information relevant, in particular, for formulating preventative and curative measures, in particular in the event of emergency health response, from importers and downstream users placing mixtures on the market. This information shall include the chemical composition of mixtures placed on the market and classified as hazardous on the basis of their health or physical effects, including the chemical identity of substances in mixtures for which a request for use of an alternative chemical name has been accepted by the Agency, in accordance with Article 24.

2. The appointed bodies shall provide all requisite guarantees for maintaining the confidentiality of the information received. Such information may only be used:

(a) to meet medical demand by formulating preventative and curative measures, in particular in the event of an emergency;

and

(b) where requested by the Member State, to undertake statistical analysis to identify where improved risk management measures may be needed.

The information shall not be used for other purposes.

3. The appointed bodies shall have at their disposal all the information required from the importers and downstream users responsible for marketing to carry out the tasks for which they are responsible.

4. By 20 January 2012 the Commission shall carry out a review to assess the possibility of harmonising the information referred to in paragraph 1, including establishing a format for the submission of information by importers and downstream users to appointed bodies. On the basis of this review, and following consultation with relevant stakeholders such as the European Association of Poison Centres and Clinical Toxicologists (EAPCCT), the Commission may adopt a Regulation adding an Annex to this Regulation.

Those measures, designed to amend non-essential elements of this Regulation, by supplementing it, shall be adopted in accordance with the regulatory procedure with scrutiny referred to in Article 54(3).

Article 46

Enforcement and reporting

1. Member States shall take all necessary measures, including maintaining a system of official controls, to ensure that substances and mixtures are not placed on the market, unless they have been classified, labelled, notified and packaged in accordance with this Regulation.

2. Member States shall submit a report to the Agency every five years by 1 July on the results of the official controls, and other enforcement measures taken. The first report shall be submitted by 20 January 2012. The Agency shall make those reports available to the Commission, which shall take them into account for its report under Article 117 of Regulation (EC) No 1907/2006.

3. The Forum referred to in Article 76(1)(f) of Regulation (EC) No 1907/2006 shall undertake the tasks specified in Article 77(4)(a) to (g) of Regulation (EC) No 1907/2006 concerning enforcement of this Regulation.

Article 47

Penalties for non-compliance

Member States shall introduce penalties for non-compliance with this Regulation and shall take all measures necessary to ensure that this Regulation is applied. The penalties must be effective, proportionate and dissuasive. Member States shall notify the Commission of the provisions for penalties by 20 June 2010 and shall notify it without delay of any subsequent amendment affecting them.

TITLE VII

COMMON AND FINAL PROVISIONS

Article 48

Advertisement

1. Any advertisement for a substance classified as hazardous shall mention the hazard classes or hazard categories concerned.

2. Any advertisement for a mixture classified as hazardous or covered by Article 25(6) which allows a member of the general public to conclude a contract for purchase without first having sight of the label shall mention the type or types of hazard indicated on the label.

The first subparagraph shall be without prejudice to Directive 97/7/EC of the European Parliament and of the Council of 20 May 1997 on the protection of consumers in respect of distance contracts ([1]).

Article 49

Obligation to maintain information and requests for information

1. The supplier shall assemble and keep available all the information used by that supplier for the purposes of classification and labelling under this Regulation for a period of at least 10 years after the substance or the mixture was last supplied by that supplier.

The supplier shall keep this information together with the information required in Article 36 of Regulation (EC) No 1907/2006.

2. In the event of a supplier ceasing activity, or transferring part or all of his operations to a third party, the party responsible for liquidating the supplier's undertaking or assuming responsibility for the placing on the market of the substance or mixture concerned shall be bound by the obligation in paragraph 1 in place of the supplier.

3. The competent authority or the enforcement authorities of a Member State in which a supplier is established or the Agency may require the supplier to submit to it any information referred to in the first subparagraph of paragraph 1.

However, where that information is available to the Agency as part of a registration pursuant to Regulation (EC) No 1907/2006 or a notification pursuant to Article 40 of this Regulation, the Agency shall use that information and the authority shall address itself to the Agency.

Article 50

Tasks of the Agency

1. The Agency shall provide the Member States and the institutions of the Community with the best possible scientific and technical advice on questions relating to chemicals which fall within its remit and which are referred to it in accordance with this Regulation.

2. The Secretariat of the Agency shall:

(a) provide industry with technical and scientific guidance and tools where appropriate on how to comply with the obligations laid down by this Regulation;

(b) provide competent authorities with technical and scientific guidance on the operation of this Regulation and provide support to the helpdesks established by Member States under Article 44.

Article 51

Free movement clause

On grounds relating to the classification, labelling or packaging of substances and mixtures within the meaning of this Regulation, Member States shall not prohibit, restrict or impede the placing on the market of substances or mixtures which comply with this Regulation and, where appropriate, with Community acts adopted in implementation of this Regulation.

Article 52

Safeguard clause

1. Where a Member State has justifiable grounds for believing that a substance or a mixture, although satisfying the requirements of this Regulation, constitutes a serious risk to human health or the environment due to reasons of classification, labelling or packaging, it may take appropriate provisional measures. The Member State shall immediately inform the Commission, the Agency and the other Member States thereof, giving the reasons for its decision.

2. Within 60 days of receipt of the information from the Member State, the Commission shall in accordance with the regulatory procedure referred to in Article 54(2) either authorise the provisional measure for a time period defined in the decision or require the Member State to revoke the provisional measure.

([1]) OJ L 144, 4.6.1997, p. 19.

3. In the case of an authorisation of a provisional measure related to classification or labelling of a substance as referred to in paragraph 2, the competent authority of the Member State concerned shall in accordance with the procedure laid down in Article 37 submit a proposal to the Agency for harmonised classification and labelling, within three months of the date of the Commission decision.

Article 53

Adaptations to technical and scientific progress

1. The Commission may adjust and adapt Articles 6(5), 11(3), 12, 14, 18(3)(b), 23, 25 to 29 and 35(2) second and third subparagraph and Annexes I to VII to technical and scientific progress, including taking due account of the further development of the GHS, in particular any UN amendments relating to the use of information on similar mixtures, and considering the developments in internationally recognised chemical programmes and of the data from accident databases. Those measures, designed to amend non-essential elements of this Regulation, shall be adopted in accordance with the regulatory procedure with scrutiny referred to in Article 54(3). On imperative grounds of urgency, the Commission may have recourse to the urgency procedure referred to in Article 54(4).

2. Member States and the Commission shall, in the manner appropriate to their role in the relevant UN fora, promote the harmonisation of the criteria for classification and labelling of persistent, bioaccumulative and toxic (PBT) and very persistent and very bioaccumulative (vPvB) substances at the level of the UN.

Article 54

Committee procedure

1. The Commission shall be assisted by the Committee instituted by Article 133 of Regulation (EC) No 1907/2006.

2. Where reference is made to this paragraph, Articles 5 and 7 of Decision 1999/468/EC shall apply, having regard to the provisions of Article 8 thereof.

The period laid down in Article 5 (6) of Decision 1999/468/EC shall be set at three months.

3. Where reference is made to this paragraph, Article 5a (1) to (4) and Article 7 of Decision 1999/468/EC shall apply, having regard to the provisions of Article 8 thereof.

4. Where reference is made to this paragraph, Article 5a(1), (2), (4) and (6) and Article 7 of Decision 1999/468/EC shall apply, having regard to the provisions of Article 8 thereof.

Article 55

Amendments to Directive 67/548/EEC

Directive 67/548/EEC shall be amended as follows:

1. in Article 1(2), the second subparagraph shall be deleted;

2. Article 4 shall be amended as follows:

(a) paragraph 3 shall be replaced by the following:

'3. Where an entry containing the harmonised classification and labelling for a particular substance has been included in Part 3 of Annex VI to Regulation (EC) No 1272/2008 of the European Parliament and of the Council of 16 December 2008 on classification, labelling and packaging of substances and mixtures (*), the substance shall be classified in accordance with that entry and paragraphs 1 and 2 shall not apply to the danger categories covered by that entry.

(*) OJ L 353, 31.12.2008, p. 1';

(b) paragraph 4 shall be deleted;

3. Article 5 shall be amended as follows:

(a) paragraph 1, second subparagraph shall be deleted;

(b) paragraph 2 shall be replaced by the following:

'2. The measures in the first subparagraph of paragraph 1 shall apply until the substance is listed in Part 3 of Annex VI to Regulation (EC) No 1272/2008 for the danger categories covered by that entry or until a decision not to list it has been taken in accordance with the procedure laid down in Article 37 of Regulation (EC) No 1272/2008.';

4. Article 6 shall be replaced by the following:

'Article 6

Obligation to carry out investigations

Manufacturers, distributors and importers of substances which appear in the EINECS but for which no entry has been included in Part 3 of Annex VI to Regulation (EC) No 1272/2008 shall carry out an investigation to make themselves aware of the relevant and accessible data which exist concerning the properties of such substances. On the basis of this information, they shall package and provisionally label dangerous substances according to the rules laid down in Articles 22 to 25 of this Directive and the criteria in Annex VI to this Directive.';

5. Article 22(3) and (4) shall be deleted;

6. Article 23(2) shall be amended as follows:

(a) in point (a), the words 'Annex I' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(b) in point (c), the words 'Annex I' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(c) in point (d), the words 'Annex I' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(d) in point (e), the words 'Annex I' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(e) in point (f), the words 'Annex I' shall be replaced by 'Part 3 of Annex VI of Regulation (EC) No 1272/2008';

7. Article 24(4) second subparagraph shall be deleted;

8. Article 28 shall be deleted;

9. Article 31(2) and (3) shall be deleted;

10. the following Article shall be inserted after Article 32:

'*Article 32a*

Transitional provision regarding labelling and packaging of substances

Articles 22 to 25 shall not apply to substances from 1 December 2010.';

11. Annex I shall be deleted.

Article 56

Amendments to Directive 1999/45/EC

Directive 1999/45/EC shall be amended as follows:

1. in Article 3(2), first indent, the words 'Annex I to Directive 67/548/EEC' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008 of the European Parliament and of the Council of 16 December 2008 on classification, labelling and packaging of substances and mixtures (*).

(*) OJ L 353, 31.12.2008, p. 1';

2. the words 'Annex I to Directive 67/548/EEC' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008' in:

(a) Article 3(3);

(b) Article 10(2), points 2.3.1, 2.3.2, 2.3.3 and 2.4 first indent;

(c) Annex II, points (a) and (b) and the last paragraph of the Introduction;

(d) Annex II, Part A,

— point 1.1.1 (a) and (b),

— point 1.2 (a) and (b),

— point 2.1.1 (a) and (b),

— point 2.2 (a) and (b),

— point 2.3 (a) and (b),

— point 3.1.1 (a) and (b),

— point 3.3 (a) and (b),

— point 3.4 (a) and (b),

— point 4.1.1 (a) and (b),

— point 4.2.1 (a) and (b),

— point 5.1.1 (a) and (b),

— point 5.2.1 (a) and (b),

— point 5.3.1 (a) and (b),

— point 5.4.1 (a) and (b),

— point 6.1 (a) and (b),

— point 6.2 (a) and (b),

— point 7.1 (a) and (b),

— point 7.2 (a) and (b),

— point 8.1 (a) and (b),

— point 8.2 (a) and (b),

— point 9.1 (a) and (b),

— point 9.2 (a) and (b),

— point 9.3 (a) and (b),

— point 9.4 (a) and (b);

(e) Annex II, the introductory paragraph of Part B;

(f) Annex III, point (a) and (b) of the Introduction;

(g) Annex III, Part A, section (a) Aquatic environment

— point 1.1 (a) and (b),

— point 2.1 (a) and (b),

— point 3.1 (a) and (b),

— point 4.1 (a) and (b),

— point 5.1 (a) and (b),

— point 6.1 (a) and (b),

(h) Annex III, Part A, section (b) Non-aquatic environment point 1.1 (a) and (b);

(i) Annex V, section A points 3 and 4;

(j) Annex V, section B point 9;

(k) Annex VI, Part A, the third column of the table under point 2;

(l) Annex VI Part B point 1, first paragraph, and the first column of the table under point 3;

(m) Annex VIII, Appendix 1, second column of the table;

(n) Annex VIII, Appendix 2, second column of the table;

3. in Annex VI, Part B, point 1, paragraph 3 first indent and paragraph 5, the words 'Annex I' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

4. in Annex VI, Part B, point 4.2, final paragraph, the words 'Annex I to Directive 67/548/EEC (19th adaptation)' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008'.

Article 57

Amendments to Regulation (EC) No 1907/2006 from the entry into force of this Regulation

Regulation (EC) No 1907/2006 shall be amended as from the entry into force of this Regulation as follows:

1. Article 14(2) shall be amended as follows:

(a) point (b) shall be replaced by the following:

'(b) the specific concentration limits that have been set in Part 3 of Annex VI to Regulation (EC) No 1272/2008 of the European Parliament and of the Council of 16 December 2008 on classification, labelling and packaging of substances and mixtures (*);

(ba) for substances classified as hazardous to the aquatic environment, if a multiplying factor (hereinafter referred to as "M-factor") has been set in Part 3 of Annex VI to Regulation (EC) No 1272/2008, the cut-off value in Table 1.1 of Annex I to that Regulation adjusted using the calculation set out in section 4.1 of Annex I to that Regulation;

(*) OJ L 353, 31.12.2008, p. 1';

(b) point (e) shall be replaced by the following:

'(e) the specific concentration limits given in an agreed entry in the classification and labelling inventory referred to in Article 42 of Regulation (EC) No 1272/2008;

(ea) for substances classified as hazardous to the aquatic environment, if an M-factor has been set in an agreed entry in the classification and labelling inventory referred to in Article 42 of Regulation (EC) No 1272/2008, the cut-off value in Table 1.1 of Annex I to that Regulation adjusted using the calculation set out in section 4.1 of Annex I to that Regulation;';

2. Article 31 shall be amended as follows:

(a) paragraph 8 shall be replaced by the following:

'8. A safety data sheet shall be provided free of charge on paper or electronically no later than the date on which the substance or mixture is first supplied.';

(b) the following paragraph shall be added:

'10. Where substances are classified in accordance with Regulation (EC) No 1272/2008 during the period from its entry into force until 1 December 2010, that classification may be added in the safety data sheet together with the classification in accordance with Directive 67/548/EEC.

From 1 December 2010 until 1 June 2015, the safety data sheets for substances shall contain the classification according to both Directive 67/548/EEC and Regulation (EC) No 1272/2008.

Where mixtures are classified in accordance with Regulation (EC) No 1272/2008 during the period from its entry into force until 1 June 2015, that classification may be added in the safety data sheet, together with the classification in accordance with Directive 1999/45/EC. However, until 1 June 2015, where substances or mixtures are both classified and labelled in accordance with Regulation (EC) No 1272/2008 that classification shall be provided in the safety data sheet, together with the classification in accordance with Directives 67/548/EEC and 1999/45/EC respectively, for the substance, the mixture and its constituents.';

3. Article 56(6)(b) shall be replaced by the following:

'(b) for all other substances, below the lowest of the concentration limits specified in Directive 1999/45/EC or in Part 3 of Annex VI to Regulation (EC) No 1272/2008 which result in the classification of the mixture as dangerous.';

4. Article 59(2) and 3 shall be amended as follows:

(a) in paragraph 2, the second sentence shall be replaced by the following:

'The dossier may be limited, if appropriate, to a reference to an entry in Part 3 of Annex VI to Regulation (EC) No 1272/2008.';

(b) in paragraph 3, the second sentence shall be replaced by the following:

'The dossier may be limited, if appropriate, to a reference to an entry in Part 3 of Annex VI to Regulation (EC) No 1272/2008.';

5. in Article 76(1)(c), the words 'Title XI' shall be replaced by 'Title V of Regulation (EC) No 1272/2008';

6. Article 77 shall be amended as follows:

(a) in paragraph 2, the first sentence of point (e) shall be replaced by the following:

'(e) establishing and maintaining database(s) with information on all registered substances, the classification and labelling inventory and the harmonised classification and labelling list established in accordance with Regulation (EC) No 1272/2008;';

(b) in paragraph 3, point (a), the words 'Titles VI to XI' shall be replaced by 'Titles VI to X';

7. Title XI shall be deleted;

8. Annex XV, sections I and II shall be amended as follows:

(a) section I shall be amended as follows:

(i) the first indent shall be deleted;

(ii) the second indent shall be replaced by the following:

'— the identification of CMRs, PBTs, vPvBs, or a substance of equivalent concern in accordance with Article 59,';

(b) in section II, point 1 shall be deleted;

9. the table in Annex XVII shall be amended as follows:

(a) the column 'Designation of the substance, of the groups of substances or of the preparation', shall be amended as follows:

(i) entries 28, 29 and 30 shall be replaced by the following:

'28. Substances which appear in Part 3 of Annex VI to Regulation (EC) No 1272/2008 classified as carcinogen category 1A or 1B (Table 3.1) or carcinogen category 1 or 2 (Table 3.2) and listed as follows:

— Carcinogen category 1A (Table 3.1)/ carcinogen category 1 (Table 3.2) listed in Appendix 1

— Carcinogen category 1B (Table 3.1)/ carcinogen category 2 (Table 3.2) listed in Appendix 2

29. Substances which appear in Part 3 of Annex VI to Regulation (EC) No 1272/2008 classified as germ cell mutagen category 1A or 1B (Table 3.1) or mutagen category 1 or 2 (Table 3.2) and listed as follows:

— Mutagen category 1A (Table 3.1)/ mutagen category 1 (Table 3.2) listed in Appendix 3

— Mutagen category 1B (Table 3.1)/ mutagen category 2 (Table 3.2) listed in Appendix 4

30. Substances which appear in Part 3 of Annex VI to Regulation (EC) No 1272/2008 classified as toxic to reproduction category 1A or 1B (Table 3.1) or toxic to reproduction category 1 or 2 (Table 3.2) and listed as follows:

— Reproductive toxicant category 1A adverse effects on sexual function and fertility or on development (Table 3.1) or reproductive toxicant category 1 with R60 (May impair fertility) or R61 (May cause harm to the unborn child) (Table 3.2) listed in Appendix 5

— Reproductive toxicant category 1B adverse effects on sexual function and fertility or on development (Table 3.1) or reproductive toxicant category 2 with R60 (May impair fertility) or R61 (May cause harm to the unborn child) (Table 3.2) listed in Appendix 6';

(b) in the column 'Conditions of restriction', in entry 28, the first indent of point 1 shall be replaced by the following:

'— either the relevant specific concentration limit specified in Part 3 of Annex VI to Regulation (EC) No 1272/2008, or';

10. Appendices 1 to 6 to Annex XVII shall be amended as follows:

(a) the Foreword shall be amended as follows:

(i) in the section entitled 'Substances', the words 'Annex I to Directive 67/548/EEC' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(ii) in the section entitled 'Index number', the words 'Annex I to Directive 67/548/EEC' shall be replaced by 'Part 3 of Annex VI to Regulation (EC) No 1272/2008';

(iii) in the section entitled 'Notes', the words 'the foreword of Annex I to Directive 67/548/EEC' shall be replaced by 'Part 1 of Annex VI to Regulation (EC) No 1272/2008';

(iv) Note A shall be replaced by the following:

'Note A:

Without prejudice to Article 17(2) of Regulation (EC) No 1272/2008, the name of the substance must appear on the label in the form of one of the designations given in Part 3 of Annex VI to that Regulation.

In that Part, use is sometimes made of a general description such as "... compounds" or "... salts". In this case, the supplier who places such a substance on the market is required to state on the label the correct name, due account being taken of Section 1.1.1.4 of Annex VI to Regulation (EC) No 1272/2008.

In accordance with Regulation (EC) No 1272/2008, where a substance is included in Part 3 of Annex VI to that Regulation, the labelling elements relevant for each specific classification covered by the entry in that Part shall be included in the label, together with the applicable label elements for any other classification not covered by that entry, and any other applicable label elements in accordance with Article 17 of that Regulation.

For substances belonging to one particular group of substances included in Part 3 of Annex VI to Regulation (EC) No 1272/2008, the labelling elements relevant for each specific classification covered by the entry in that Part shall be included in the label, together with the applicable label elements for any other classification not covered by that entry, and any other applicable label elements in accordance with Article 17 of that Regulation.

For substances belonging to more than one group of substances included in Part 3 of Annex VI to Regulation (EC) No 1272/2008, the labelling elements relevant for each specific classification covered by both entries in that Part shall be included in the label, together with the applicable label elements for any other classification not covered by that entry, and any other applicable label elements in accordance with Article 17 of that Regulation. In cases where two different classifications are given in the two entries for the same hazard class or differentiation, the classification reflecting the more severe classification shall be used.';

(v) Note D shall be replaced by the following:

'Note D:

Certain substances which are susceptible to spontaneous polymerisation or decomposition are generally placed on the market in a stabilised form. It is in this form that they are listed in Part 3 of Annex VI to Regulation (EC) No 1272/2008.

However, such substances are sometimes placed on the market in a non-stabilised form. In this case, the supplier who places such a substance on the market must state on the label the name of the substance followed by the words "non-stabilised".';

(vi) Note E shall be deleted;

(vii) Note H shall be replaced by the following:

'Note H:

The classification and label shown for this substance applies to the hazard or hazards indicated by the hazard statement or hazard statements in combination with the hazard classification shown. The requirements of Article 4 of Regulation (EC) No 1272/2008 on suppliers of this substance apply to all other hazard classes, differentiations and categories.

The final label shall follow the requirements of section 1.2 of Annex I to Regulation (EC) No 1272/2008.';

(viii) Note K shall be replaced by the following:

'Note K:

The classification as a carcinogen or mutagen need not apply if it can be shown that the substance contains less than 0,1 % w/w 1,3-butadiene (Einecs No 203-450-8). If the substance is not classified as a carcinogen or mutagen, at least the precautionary statements (P102-)P210-P403 should apply. This note applies only to certain complex oil-derived substances in Part 3 of Annex VI to Regulation (EC) No 1272/2008.';

(ix) Note S shall be replaced by the following:

'Note S:

This substance may not require a label according to Article 17 of Regulation (EC) No 1272/2008 (see section 1.3 of Annex I to that Regulation).';

(b) in Appendix 1, the title shall be replaced by the following:

'Point 28 — Carcinogens: category 1A (Table 3.1)/ category 1 (Table 3.2)';

(c) Appendix 2 shall be amended as follows:

(i) the title shall be replaced by 'Point 28 — Carcinogens: category 1B (Table 3.1)/ category 2 (Table 3.2)';

(ii) in the entries index Nos 024-017-00-8, 611-024-001, 611-029-00-9, 611-030-00-4 and 650-017-00-8, the words 'Annex I to Directive 67/548/EEC' shall be replaced by 'Annex VI to Regulation (EC) No 1272/2008.';

(d) in Appendix 3, the title shall be replaced by the following:

'Point 29 — Mutagens: category 1A (Table 3.1)/ category 1 (Table 3.2)';

(e) in Appendix 4, the title shall be replaced by the following:

'Point 29 — Mutagens: category 1B (Table 3.1)/ category 2 (Table 3.2)';

(f) in Appendix 5, the title shall be replaced by the following:

'Point 30 — Reproductive toxicants: category 1A (Table 3.1)/category 1 (Table 3.2)';

(g) in Appendix 6, the title shall be replaced by the following:

'Point 30 — Reproductive toxicants: category 1B (Table 3.1)/category 2 (Table 3.2)';

11. the word 'preparation' or 'preparations' within the meaning of Article 3 (2) of Regulation (EC) 1907/2006 shall be replaced by 'mixture' or 'mixtures' respectively throughout the text.

Article 58

Amendments to Regulation (EC) No 1907/2006 from 1 December 2010

Regulation (EC) No 1907/2006 shall be amended from 1 December 2010 as follows:

1. in Article 14(4), the introductory sentence shall be replaced by the following:

'4. If, as a result of carrying out steps (a) to (d) of paragraph 3, the registrant concludes that the substance fulfils the criteria for any of the following hazard classes or categories set out in Annex I to Regulation (EC) No 1272/2008:

(a) hazard classes 2.1 to 2.4, 2.6 and 2.7, 2.8 types A and B, 2.9, 2.10, 2.12, 2.13 categories 1 and 2, 2.14 categories 1 and 2, 2.15 types A to F;

(b) hazard classes 3.1 to 3.6, 3.7 adverse effects on sexual function and fertility or on development, 3.8 effects other than narcotic effects, 3.9 and 3.10;

(c) hazard class 4.1;

(d) hazard class 5.1,

or is assessed to be a PBT or vPvB, the chemical safety assessment shall include the following additional steps:';

2. Article 31 shall be amended as follows

(a) paragraph 1(a) shall be replaced by the following:

'(a) where a substance meets the criteria for classification as hazardous in accordance with Regulation (EC) No 1272/2008 or a mixture meets the criteria for classification as dangerous in accordance with Directive 1999/45/EC; or';

(b) paragraph 4 shall be replaced by the following:

'4. The safety data sheet need not be supplied where substances that are hazardous in accordance with Regulation (EC) No 1272/2008 or mixtures that are dangerous in accordance with Directive 1999/45/EC, offered or sold to the general public, are provided with sufficient information to enable users to take the necessary measures as regards the protection of human health, safety and the environment, unless requested by a downstream user or distributor.';

3. Article 40(1) shall be replaced by the following:

'1. The Agency shall examine any testing proposal set out in a registration or a downstream user report for provision of the information specified in Annexes IX and X for a substance. Priority shall be given to registrations of substances which have or may have PBT, vPvB, sensitising and/or carcinogenic, mutagenic or toxic for reproduction (CMR) properties, or substances above 100 tonnes per year with uses resulting in widespread and diffuse exposure, provided they fulfil the criteria for any of the following hazard classes or categories set out in Annex I of Regulation (EC) No 1272/2008:

(a) hazard classes 2.1 to 2.4, 2.6 and 2.7, 2.8 types A and B, 2.9, 2.10, 2.12, 2.13 categories 1 and 2, 2.14 categories 1 and 2, 2.15 types A to F;

(b) hazard classes 3.1 to 3.6, 3.7 adverse effects on sexual function and fertility or on development, 3.8 effects other than narcotic effects, 3.9 and 3.10;

(c) hazard class 4.1;

(d) hazard class 5.1.';

4. Article 57(a), (b) and (c) shall be replaced by the following:

'(a) substances meeting the criteria for classification in the hazard class carcinogenicity category 1A or 1B in accordance with section 3.6 of Annex I to Regulation (EC) No 1272/2008;

(b) substances meeting the criteria for classification in the hazard class germ cell mutagenicity category 1A or 1B in accordance with section 3.5 of Annex I to Regulation (EC) No 1272/2008;

(c) substances meeting the criteria for classification in the hazard class reproductive toxicity category 1A or 1B, adverse effects on sexual function and fertility or on development in accordance with section 3.7 of Annex I to Regulation(EC) No 1272/2008;';

5. in Article 65 the words 'Directive 67/548/EEC' shall be replaced by 'Directive 67/548/EEC and Regulation (EC) No 1272/2008';

6. Article 68(2) shall be replaced by the following:

'2. For a substance on its own, in a mixture or in an article which meets the criteria for classification in the hazard classes carcinogenicity, germ cell mutagenicity or reproductive toxicity, category 1A or 1B, and could be used by consumers and for which restrictions to consumer use are proposed by the Commission, Annex XVII shall be amended in accordance with the procedure referred to in Article 133(4). Articles 69 to 73 shall not apply.';

7. Article 119 shall be amended as follows:

(a) in paragraph 1, point (a) shall be replaced by the following:

'(a) without prejudice to paragraph 2(f) and (g) of this Article, the name in the IUPAC nomenclature for substances fulfilling the criteria for any of the following hazard classes or categories set out in Annex I to Regulation (EC) No 1272/2008:

— hazard classes 2.1 to 2.4, 2.6 and 2.7, 2.8 types A and B, 2.9, 2.10, 2.12, 2.13 categories 1 and 2, 2.14 categories 1 and 2, 2.15 types A to F;

— hazard classes 3.1 to 3.6, 3.7 adverse effects on sexual function and fertility or on development, 3.8 effects other than narcotic effects, 3.9 and 3.10;

— hazard class 4.1;

— hazard class 5.1.';

(b) paragraph 2 shall be amended as follows:

(i) point (f) shall be replaced by the following:

'(f) subject to Article 24 of Regulation (EC) No 1272/2008, the name in the IUPAC nomenclature for non-phase-in substances referred to in paragraph 1(a) of this Article for a period of six years;'

(ii) in point (g), the introductory phrase shall be replaced by the following:

'(g) subject to Article 24 of Regulation (EC) No 1272/2008, the name in the IUPAC nomenclature for substances referred to in paragraph 1(a) of this Article that are only used as one or more of the following:';

8. in Article 138(1), the second sentence of the introductory phrase shall be replaced by the following:

'However, for substances meeting the criteria for classification in the hazard classes carcinogenicity, germ cell mutagenicity or reproductive toxicity, category 1A or 1B, in accordance with Regulation (EC) No 1272/2008, the review shall be carried out by 1 June 2014.';

9. Annex III shall be amended as follows:

(a) point (a) shall be replaced by the following:

'(a) substances for which it is predicted (i.e. by the application of (Q)SARs or other evidence) that they are likely to meet the criteria for category 1A or 1B classification in the hazard classes carcinogenicity, germ cell mutagenicity or reproductive toxicity or the criteria in Annex XIII;';

(b) in point (b), point (ii) shall be replaced by the following:

'(ii) for which it is predicted (i.e. by application of (Q) SARs or other evidence) that they are likely to meet the classification criteria for any health or environmental hazard classes or differentiations under Regulation (EC) No 1272/2008.';

10. in Annex V, point 8, the words 'Directive 67/548/EEC' shall be replaced by 'Regulation (EC) No 1272/2008';

11. in Annex VI, sections 4.1, 4.2 and 4.3 shall be replaced by the following:

'4.1 The hazard classification of the substance(s), resulting from the application of Title I and II of Regulation (EC) No 1272/2008 for all hazard classes and categories in that Regulation,

In addition, for each entry, the reasons why no classification is given for a hazard class or differentiation of a hazard class should be provided (i.e. if data are lacking, inconclusive, or conclusive but not sufficient for classification),

4.2 The resulting hazard label for the substance(s), resulting from the application of Title III of Regulation (EC) No 1272/2008,

4.3 Specific concentration limits, where applicable, resulting from the application of Article 10 of Regulation (EC) No 1272/2008 and Articles 4 to 7 of Directive 1999/45/EC.';

12. Annex VIII shall be amended as follows:

(a) in column 2, the second indent of point 8.4.2 shall be replaced by the following:

'— the substance is known to be carcinogenic category 1A or 1B or germ cell mutagenic category 1A, 1B or 2.';

(b) in column 2, the second and third paragraphs of point 8.7.1 shall be replaced by the following:

'If a substance is known to have an adverse effect on fertility, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage fertility (H360F), and the available data are adequate to support a robust risk assessment, then no further testing for fertility will be necessary. However, testing for developmental toxicity must be considered.

If a substance is known to cause developmental toxicity, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage the unborn child (H360D), and the available data are adequate to support a robust risk assessment, then no further testing for developmental toxicity will be necessary. However, testing for effects on fertility must be considered.';

13. in Annex IX, column 2, point 8.7, the second and third paragraphs shall be replaced by the following:

'If a substance is known to have an adverse effect on fertility, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage fertility (H360F), and the available data are adequate to support a robust risk assessment, then no further testing for fertility will be necessary. However, testing for developmental toxicity must be considered.

If a substance is known to cause developmental toxicity, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage the unborn child (H360D), and the available data are adequate to support a robust risk assessment, then no further testing for developmental toxicity will be necessary. However, testing for effects on fertility must be considered.';

14. Annex X shall be amended as follows:

(a) in column 2, point 8.7, the second and third paragraphs shall be replaced by the following:

'If a substance is known to have an adverse effect on fertility, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage fertility (H360F), and the available data are adequate to support a robust risk assessment, then no further testing for fertility will be necessary. However, testing for developmental toxicity must be considered.

If a substance is known to cause developmental toxicity, meeting the criteria for classification as toxic for reproduction category 1A or 1B: May damage the unborn child (H360D), and the available data are adequate to support a robust risk assessment, then no further testing for developmental toxicity will be necessary. However, testing for effects on fertility must be considered.'

(b) in column 2, point 8.9.1, the second indent of the first paragraph shall be replaced by the following:

'— the substance is classified as germ cell mutagen category 2 or there is evidence from the repeated dose study(ies) that the substance is able to induce hyperplasia and/or pre-neoplastic lesions.'

(c) in column 2, the second paragraph of point 8.9.1 shall be replaced by the following:

'If the substance is classified as germ cell mutagen category 1A or 1B, the default presumption would be that a genotoxic mechanism for carcinogenicity is likely. In these cases, a carcinogenicity test will normally not be required.';

15. in Annex XIII, the second and third indents of point 1.3 shall be replaced by the following:

'— the substance is classified as carcinogenic (category 1A or 1B), germ cell mutagenic (category 1A or 1B), or toxic for reproduction (category 1A, 1B or 2), or

— there is other evidence of chronic toxicity, as identified by the classifications STOT (repeated exposure), category 1 (oral, dermal, inhalation of gases/vapours, inhalation of dust/mist/fume) or category 2 (oral, dermal, inhalation of gases/vapours, inhalation of dust/mist/fume) according to Regulation (EC) No 1272/2008';

16. in the table in Annex XVII, the column 'Designation of the substance, of the groups of substances or of the mixture' shall be amended as follows:

(a) entry 3 shall be replaced by the following:

'3. Liquid substances or mixtures which are regarded as dangerous in accordance with Directive 1999/45/EC or are fulfilling the criteria for any of the following hazard classes or categories set out in Annex I to Regulation (EC) No 1272/2008:

(a) hazard classes 2.1 to 2.4, 2.6 and 2.7, 2.8 types A and B, 2.9, 2.10, 2.12, 2.13 categories 1 and 2, 2.14 categories 1 and 2, 2.15 types A to F;

(b) hazard classes 3.1 to 3.6, 3.7 adverse effects on sexual function and fertility or on development, 3.8 effects other than narcotic effects, 3.9 and 3.10;

(c) hazard class 4.1;

(d) hazard class 5.1.';

(b) entry 40 shall be replaced by the following:

'40. Substances classified as flammable gases category 1 or 2, flammable liquids categories 1, 2 or 3, flammable solids category 1 or 2, substances and mixtures which, in contact with water, emit flammable gases, category 1, 2 or 3, pyrophoric liquids category 1 or pyrophoric solids category 1, regardless of whether they appear in Part 3 of Annex VI to that Regulation or not.'.

Article 59

Amendments to Regulation (EC) No 1907/2006 from 1 June 2015

Regulation (EC) No 1907/2006 shall be amended from 1 June 2015 as follows:

1. Article 14(2) shall be replaced by the following:

'2. A chemical safety assessment in accordance with paragraph 1 need not be performed for a substance which is present in a mixture if the concentration of the substance in the mixture is less than

(a) the cut-off value referred to in Article 11, paragraph 3 of Regulation (EC) No 1272/2008;

(b) 0,1 % weight by weight (w/w), if the substance meets the criteria in Annex XIII to this Regulation.';

2. Article 31 shall be amended as follows:

(a) in paragraph 1, point (a) shall be replaced by the following:

'(a) where a substance or mixture meets the criteria for classification as hazardous in accordance with Regulation (EC) No 1272/2008; or';

(b) paragraph 3 shall be replaced by the following:

'3. The supplier shall provide the recipient at his request with a safety data sheet compiled in accordance with Annex II, where a mixture does not meet the criteria for classification as hazardous in accordance with Titles I and II of Regulation (EC) No 1272/2008, but contains:

(a) in an individual concentration of ≥ 1 % by weight for non-gaseous mixtures and ≥ 0,2 % by volume for gaseous mixtures at least one substance posing human health or environmental hazards; or

(b) in an individual concentration of ≥ 0,1 % by weight for non-gaseous mixtures at least one substance that is carcinogenic category 2 or toxic to reproduction category 1A, 1B and 2, skin sensitiser category 1, respiratory sensitiser category 1, or has effects on or via lactation or is persistent, bioaccumulative and toxic (PBT) in accordance with the criteria set out in Annex XIII or very persistent and very bioaccumulative (vPvB) in accordance with the criteria set out in Annex XIII or has been included for reasons other than those referred to in point (a) in the list established in accordance with Article 59(1); or

(c) a substance for which there are Community workplace exposure limits';

(c) paragraph 4 shall be replaced by the following:

'4. The safety data sheet need not be supplied where hazardous substances or mixtures offered or sold to the general public are provided with sufficient information to enable users to take the necessary measures as regards the protection of human health, safety and the environment, unless requested by a downstream user or distributor.';

3. Article 56(6)(b) shall be replaced by the following:

'(b) for all other substances, below the values specified in Article 11(3) of Regulation (EC) No 1272/2008 which result in the classification of the mixture as hazardous.';

4. in Article 65 the words 'and Directive 1999/45/EC' shall be deleted;

5. Annex II shall be amended as follows:

(a) point 1.1 shall be replaced by:

'1.1. Identification of the substance or mixture

The term used for identification of a substance shall be identical to that provided on the label in accordance with Article 18(2) of Regulation (EC) No 1272/2008.

The term used for identification of a mixture shall be identical to that provided on the label in accordance with Article 18(3)(a) of Regulation (EC) No 1272/2008.';

(b) footnote 1 to point 3.3(a), first indent, shall be deleted;

(c) point 3.6 shall be replaced by:

'3.6. Where, in accordance with Article 24 of Regulation (EC) No 1272/2008, the Agency has agreed that the chemical identity of a substance may be kept confidential on the label and in the safety data sheet, their chemical nature shall be described under heading 3 in order to ensure safe handling.

The name used on the safety data sheet (including for the purposes of paragraphs 1.1, 3.2, 3.3 and 3.5) shall be the same as that used on the label, agreed in accordance with the procedure set out in Article 24 of Regulation (EC) No 1272/2008.';

6. in Annex VI section 4.3 shall be replaced by the following:

'4.3 Specific concentration limits, where applicable, resulting from the application of Article 10 of Regulation (EC) No 1272/2008.';

7. Annex XVII shall be amended as follows:

(a) in the column 'Designation of the substance, of the groups of substances or of the mixture' of the table in entry 3, the words 'which are regarded as dangerous in accordance with Directive 1999/45/EC or are' shall be deleted;

(b) in the column 'Conditions of restriction' of the table, entry 28 shall be amended as follows:

(i) the second indent of point 1 shall be replaced by the following:

'— the relevant generic concentration limit specified in Part 3 of Annex I of Regulation (EC) No 1272/2008.';

(ii) point 2 (d) shall be replaced by the following:

'(d) artists' paints covered by Regulation (EC) No 1272/2008.'.

Article 60

Repeal

Directive 67/548/EEC and Directive 1999/45/EC shall be repealed with effect from 1 June 2015.

Article 61

Transitional provisions

1. Until 1 December 2010, substances shall be classified, labelled and packaged in accordance with Directive 67/548/EEC.

Until 1 June 2015, mixtures shall be classified, labelled and packaged in accordance with Directive 1999/45/EC.

2. By way of derogation from the second subparagraph of Article 62 of this Regulation and in addition to the requirements of paragraph 1 of this Article, substances and mixtures may, before 1 December 2010 and 1 June 2015 respectively, be classified, labelled and packaged in accordance with this Regulation. In that case, the provisions on labelling and packaging in Directives 67/548/EEC and 1999/45/EC shall not apply.

3. From 1 December 2010 until 1 June 2015, substances shall be classified in accordance with both Directive 67/548/EEC and this Regulation. They shall be labelled and packaged in accordance with this Regulation.

4. By way of derogation from the second subparagraph of Article 62 of this Regulation, substances classified, labelled and packaged in accordance with Directive 67/548/EEC and already placed on the market before 1 December 2010, are not required to be relabelled and repackaged in accordance with this Regulation until 1 December 2012.

By way of derogation from the second subparagraph of Article 62 of this Regulation, mixtures classified, labelled and packaged in accordance with Directive 1999/45/EC and already placed on the market before 1 June 2015 are not required to be relabelled and repackaged in accordance with this Regulation until 1 June 2017.

5. Where a substance or mixture has been classified in accordance with Directive 67/548/EEC or 1999/45/EC before 1 December 2010 or 1 June 2015 respectively, manufacturers, importers and downstream users may amend the classification of the substance or mixture using the conversion table in Annex VII to this Regulation.

6. Until 1 December 2011 a Member State may maintain any existing and more stringent classification and labelling of substances entered into Part 3 of Annex VI to this Regulation, provided that these classifications and labelling elements have been notified to the Commission in accordance with the safeguard clause in Directive 67/548/EEC before 20 January 2009 and that the Member State submits a proposal for harmonised classification and labelling containing these classifications and labelling elements to the Agency in accordance with Article 37(1) of this Regulation by 1 June 2009.

It is a precondition that a decision on the proposed classification and labelling by the Commission in accordance with the safeguard clause of Directive 67/548/EEC has not yet been taken before 20 January 2009.

If the proposed harmonised classification and labelling submitted under the first subparagraph is not included or is included in an amended form in Part 3 of Annex VI in accordance with Article 37(5), the exemption in the first subparagraph of this paragraph is no longer valid.

Article 62

Entry into force

This Regulation shall enter into force on the 20th day following its publication in the *Official Journal of the European Union*.

Titles II, III and IV shall apply in respect of substances from 1 December 2010 and in respect of mixtures from 1 June 2015.

This Regulation shall be binding in its entirety and directly applicable in all Member States.

Done at Strasbourg, 16 December 2008.

For the European Parliament	*For the Council*
The President	*The President*
H.-G. PÖTTERING	B. LE MAIRE

ANNEX I

CLASSIFICATION AND LABELLING REQUIREMENTS FOR HAZARDOUS SUBSTANCES AND MIXTURES

This annex sets out the criteria for classification in hazard classes and in their differentiations and sets out additional provisions on how the criteria may be met.

1. PART 1: GENERAL PRINCIPLES FOR CLASSIFICATION AND LABELLING

1.0. Definitions

Gas means a substance which:

(i) at 50 °C has a vapour pressure greater than 300 kPa (absolute); or

(ii) is completely gaseous at 20 °C at a standard pressure of 101,3 kPa;

Liquid means a substance or mixture which:

(i) at 50 °C has a vapour pressure of not more than 300 kPa (3 bar);

(ii) is not completely gaseous at 20 °C and at a standard pressure of 101,3 kPa; and

(iii) which has a melting point or initial melting point of 20 °C or less at a standard pressure of 101,3 kPa;

Solid means a substance or mixture which does not meet the definitions of liquid or gas.

1.1. Classification of substances and mixtures

1.1.0. *Cooperation to meet the requirements in this Regulation*

Suppliers in a supply chain shall cooperate to meet the requirements for classification, labelling and packaging set out in this Regulation.

Suppliers in an industry sector may cooperate to manage the transitional arrangements in Article 61 for substances and mixtures placed on the market.

Suppliers in an industry sector may cooperate through formation of a network or by other means to share data and expertise when classifying substances and mixtures in accordance with Title II of this Regulation. In these circumstances suppliers in an industry sector shall document fully the basis on which classification decisions are made and shall make available to the competent authorities and, on request, to the relevant enforcement authorities the documentation, together with the data and information on which classifications are based. However, where suppliers in an industry sector cooperate in this way, each supplier shall remain fully responsible for the classification, labelling and packaging of substances and mixtures he places on the market, and for meeting any other requirements of this Regulation.

The network may also be used to exchange information and best practices with a view to simplifying fulfilment of the notification obligations.

1.1.1. *The role and application of expert judgement and weight of evidence determination*

1.1.1.1. Where the criteria cannot be applied directly to available identified information, or where only the information referred to in Article 6(5) is available, the weight of evidence determination using expert judgment shall be applied in accordance with Article 9(3) or 9(4) respectively.

1.1.1.2. The approach to classifying mixtures may include the application of expert judgement in a number of areas in order to ensure existing information can be used for as many mixtures as possible in order to provide protection for human health and the environment. Expert judgement may also be required in interpreting data for hazard classification of substances, especially where weight of evidence determinations are needed.

1.1.1.3. A weight of evidence determination means that all available information bearing on the determination of hazard is considered together, such as the results of suitable in vitro tests, relevant animal data, information from the application of the category approach (grouping, read-across), (Q)SAR results, human experience such as occupational data and data from accident databases, epidemiological and clinical studies and well-documented case reports and observations. The quality and consistency of the data shall be given appropriate weight. Information on substances or mixtures related to the substance or mixture being classified shall be considered as appropriate, as well as site of action and mechanism or mode of action study results. Both positive and negative results shall be assembled together in a single weight of evidence determination.

1.1.1.4. For the purpose of classification for health hazards (Part 3) established hazardous effects seen in appropriate animal studies or from human experience that are consistent with the criteria for classification shall normally justify classification. Where evidence is available from both humans and animals and there is a conflict between the findings, the quality and reliability of the evidence from both sources shall be evaluated in order to resolve the question of classification. Generally, adequate, reliable and representative data on humans (including epidemiological studies, scientifically valid case studies as specified in this Annex or statistically backed experience) shall have precedence over other data. However, even well-designed and conducted epidemiological studies may lack a sufficient number of subjects to detect relatively rare but still significant effects, to assess potentially confounding factors. Therefore, positive results from well-conducted animal studies are not necessarily negated by the lack of positive human experience but require an assessment of the robustness, quality and statistical power of both the human and animal data.

1.1.1.5. For the purpose of classification for health hazards (Part 3) route of exposure, mechanistic information and metabolism studies are pertinent to determining the relevance of an effect in humans. When such information, as far as there is reassurance about the robustness and quality of the data, raises doubt about relevance in humans, a lower classification may be warranted. When there is scientific evidence that the mechanism or mode of action is not relevant to humans, the substance or mixture should not be classified.

1.1.2. ***Specific concentration limits, M-factors and generic cut-off values***

1.1.2.1. Specific concentration limits or M-factors shall be applied in accordance with Article 10.

1.1.2.2. *Cut-off values*

1.1.2.2.1. Cut-off values indicate when the presence of a substance needs to be taken into account for the purposes of classification of a substance or a mixture containing that hazardous substance, whether as an identified impurity, additive, or individual constituent (see Article 11).

1.1.2.2.2. The cut-off values referred to in Article 11 shall be the following:

(a) For health and environmental hazards in Parts 3, 4 and 5 of this Annex:

(i) for substances where a specific concentration limit is set for the relevant hazard class or differentiation either in Part 3 of Annex VI or in the classification and labelling inventory referred to in Article 42, and where the hazard class or differentiation is mentioned in Table 1.1, the lower of the specific concentration limit and the relevant generic cut-off value in Table 1.1; or

(ii) for substances where a specific concentration limit is set for the relevant hazard class or differentiation either in Part 3 of Annex VI or in the classification and labelling inventory referred to in Article 42, and where the hazard class or differentiation is not mentioned in Table 1.1, the specific concentration limit set either in Part 3 of Annex VI or in the classification and labelling inventory; or

(iii) for substances where no specific concentration limit is set for the relevant hazard class or differentiation either in Part 3 of Annex VI or in the classification and labelling inventory referred to in Article 42, and where the hazard class or differentiation is mentioned in Table 1.1, the relevant generic cut-off value set out in that table; or

(iv) for substances where no specific concentration limit is set for the relevant hazard class or differentiation either in Part 3 of Annex VI or in the classification and labelling inventory referred to in Article 42, and where the hazard class or differentiation is not mentioned in Table 1.1, the generic concentration limit for classification in the relevant sections of Parts 3, 4 and 5 of this Annex.

(b) For aquatic environmental hazards in section 4.1 of this Annex:

(i) for substances where an M-factor has been set for the relevant hazard category either in Part 3 of Annex VI, or in the classification and labelling inventory referred to in Article 42, the generic cut-off value in Table 1.1 adjusted using the calculation set out in section 4.1 of this Annex; or

(ii) for substances where no M-factor is set for the relevant hazard category either in Part 3 of Annex VI or in the classification and labelling inventory referred to in Article 42, the relevant generic cut-off value set out in Table 1.1.

Table 1.1

Generic cut-off values

Hazard class	Generic cut-off values to be taken into account
Acute Toxicity:	
— Category 1-3	0,1 %
— Category 4	1 %
Skin corrosion/Irritation	1 % ([1])
Serious damage to eyes/eye irritation	1 % ([2])
Hazardous to Aquatic Environment	
— Acute Category 1	0,1 % ([3])
— Chronic Category 1	0,1 % ([3])
— Chronic Category 2-4	1 %

([1]) Or < 1 % where relevant, see 3.2.3.3.1.
([2]) Or < 1 % where relevant, see 3.3.3.3.1.
([3]) Or < 0,1 % where relevant, see 4.1.3.1.

Note

Generic cut-off values are in weight percentages except for gaseous mixtures where they are in volume percentage.

1.1.3. ***Bridging principles for the classification of mixtures where test data are not available for the complete mixture***

Where the mixture itself has not been tested to determine its hazardous properties, but there are sufficient data on similar tested mixtures and individual hazardous ingredient substances to adequately characterise the hazards of the mixture, these data shall be used in accordance with the following bridging rules referred to in Article 9(4) for each individual hazard class in Part 3 and Part 4 of this Annex, subject to any specific provisions for mixtures in each hazard class.

1.1.3.1. *Dilution*

If a mixture is diluted with a substance (diluent) which has an equivalent or lower hazard category classification than the least hazardous original ingredient substance and which is not expected to affect the hazard classification of other ingredient substances, then one of the following shall be applied:

— the new mixture shall be classified as equivalent to the original mixture;

— the method explained in each section of Part 3 and in Part 4 for classification of mixtures when data are available for all components or only some components of the mixture;

— in the case of acute toxicity, the method for classification of mixtures based on ingredients of the mixture (additivity formula).

1.1.3.2. *Batching*

The hazard category of one production batch of a mixture can be assumed to be substantially equivalent to that of another production batch of the same commercial product, and produced by or under the control of the same supplier, unless there is reason to believe there is significant variation such that the hazard classification of the batch has changed. If the latter occurs, a new evaluation is necessary.

1.1.3.3. *Concentration of highly hazardous mixtures*

In the case of the classification of mixtures covered by sections 3.1, 3.2, 3.3, 3.8, 3.9, 3.10 and 4.1, if a mixture is classified in the highest hazard category or sub-category, and the concentration of the ingredients of the mixture that are in that category or sub-category is increased, the new mixture shall be classified in that category or sub-category without additional testing.

1.1.3.4. *Interpolation within one toxicity category*

In the case of the classification of mixtures covered by sections 3.1, 3.2, 3.3, 3.8, 3.9, 3.10 and 4.1, for three mixtures with identical hazardous ingredients, where mixtures A and B are in the same hazard category and mixture C has the same active hazardous ingredients with concentrations intermediate to the concentrations of those hazardous ingredients in mixtures A and B, then mixture C is assumed to be in the same hazard category as A and B.

1.1.3.5. *Substantially similar mixtures*

Given the following:

(a) two mixtures each containing two ingredients:

(i) A + B

(ii) C + B;

(b) the concentration of ingredient B is essentially the same in both mixtures;

(c) the concentration of ingredient A in mixture (i) equals that of ingredient C in mixture (ii);

(d) hazard data for A and C are available and substantially equivalent, i.e. they are in the same hazard category and are not expected to affect the hazard classification of B.

If mixture (i) is already classified in a particular hazard class based on test data, mixture (ii) shall be assigned the same hazard category.

1.1.3.6. *Review of classification where the composition of a mixture has changed*

The following variations in initial concentration are defined for the application of Article 15(2)(a):

Table 1.2

Bridging Principle for changes in the composition of a mixture

Initial concentration range of the constituent	Permitted variation in initial concentration of the constituent
≤ 2,5 %	± 30 %
2,5 < C ≤ 10 %	± 20 %
10 < C ≤ 25 %	± 10 %
25 < C ≤ 100 %	± 5 %

1.1.3.7. *Aerosols*

In the case of the classification of mixtures covered by sections 3.1, 3.2, 3.3, 3.4, 3.8 and 3.9, an aerosol form of a mixture shall be classified in the same hazard category as the non-aerosolised form of the mixture, provided that the added propellant does not affect the hazardous properties of the mixture upon spraying and scientific evidence is available demonstrating that the aerosolised form is not more hazardous than the nonaerosolised form.

1.2. **Labelling**

1.2.1. ***Dimensions and make-up of the label elements***

1.2.1.1. Hazard pictograms as laid down in Annex V shall have a black symbol on a white background with a red frame sufficiently wide to be clearly visible.

1.2.1.2. Hazard pictograms shall be in the shape of a square set at a point. Each hazard pictogram shall cover at least one fifteenth of the surface area of the harmonised label but the minimum area shall not be less than 1 cm^2.

1.2.1.3. The dimensions of the label shall be as follows:

Table 1.3

Dimension of labels

Capacity of the package	Dimensions (in millimetres)
Not exceeding 3 litres:	If possible, at least 52 × 74
Greater than 3 litres but, not exceeding 50 litres:	At least 74 × 105
Greater than 50 litres but not exceeding 500 litres:	At least 105 × 148
Greater than 500 litres:	At least 148 × 210

1.3. **Derogations from labelling requirements for special cases**

In accordance with Article 23 the following derogations shall apply:

1.3.1. ***Transportable gas cylinders***

For transportable gas cylinders, one of the following shall be permitted to be used for gas cylinders with a water capacity of less than or equal to 150 litres:

(a) A format and dimensions following the prescriptions of the current edition of Standard ISO 7225 relating to 'Gas cylinders — Precautionary labels'. In this case, the label can bear the generic name or industrial or commercial name of the substance or mixture provided that the hazardous substances in a mixture are shown on the body of the gas cylinder in a clear and indelible way.

(b) The information specified in Article 17 provided on a durable information disc or label held captive on the cylinder.

1.3.2. ***Gas containers intended for propane, butane or liquefied petroleum gas (LPG)***

1.3.2.1. If propane, butane and liquefied petroleum gas or a mixture containing these substances classified in accordance with the criteria of this Annex, is placed on the market in closed refillable cylinders or in non-refillable cartridges within the scope of EN 417 as fuel gases which are only released for combustion (current edition of EN 417, relating to 'Non-refillable metallic gas cartridges for liquefied petroleum gases, with or without a valve, for use with portable appliances; construction, inspection, testing and marking'), these cylinders or cartridges shall only be labelled with the appropriate pictogram and the hazard and precautionary statements concerning flammability.

1.3.2.2. No information concerning the effects on human health and the environment is required on the label. Instead the supplier shall provide the information concerning effects on human health and the environment to downstream users or distributors by means of the safety data sheet (SDS).

1.3.2.3. For consumers, sufficient information shall be transmitted to enable them to take all necessary measures for health and safety.

1.3.3. ***Aerosols and containers fitted with a sealed spray attachment and containing substances or mixtures classified as presenting an aspiration hazard***

With regard to the application of section 3.10.4, substances or mixtures classified in accordance with the criteria of sections 3.10.2 and 3.10.3 need not be labelled for this hazard when placed on the market in aerosol containers or in containers fitted with a sealed spray attachment.

1.3.4. ***Metals in massive form, alloys, mixtures containing polymers, mixtures containing elastomers***

1.3.4.1. Metals in massive form, alloys, mixtures containing polymers and mixtures containing elastomers do not require a label according to this Annex, if they do not present a hazard to human health by inhalation, ingestion or contact with skin or to the aquatic environment in the form in which they are placed on the market, although classified as hazardous in accordance with the criteria of this Annex.

1.3.4.2. Instead, the supplier shall provide the information to downstream users or distributors by means of the SDS.

1.3.5. ***Explosives placed on the market with a view to obtaining an explosive or pyrotechnic effect***

Explosives, as referred to in section 2.1, placed on the market with a view to obtaining an explosive or pyrotechnic effect shall be labelled and packaged in accordance with the requirements for explosives only.

1.4. **Request for use of an alternative chemical name**

1.4.1. ***Requests for use of an alternative chemical name under Article 24 may be granted only where***

(I) the substance has not been assigned a Community workplace exposure limit; and

(II) the manufacturer, importer or downstream user can demonstrate that the use of the alternative chemical name meets the need to provide enough information for necessary health and safety precautions to be taken in the workplace and the need to ensure that risks from handling the mixture can be controlled; and

(III) the substance is classified exclusively as one or more of the following hazard categories:

(a) any of the hazard categories referred to in Part 2 of this Annex;

(b) Acute toxicity, Category 4;

(c) Skin corrosion/irritation, Category 2;

(d) Serious eye damage/eye irritation, Category 2;

(e) Specific target organ toxicity — Single exposure, Category 2 or 3;

(f) Specific target organ toxicity — Repeated exposure, Category 2;

(g) Hazardous to the aquatic environment — Chronic, Category 3 or 4.

1.4.2. ***The choice of the chemical name(s) for mixtures intended for the fragrance or perfume industry***

In the case of substances occurring in nature, a chemical name or chemical names of the type 'essential oil of …' or 'extract of …' may be used instead of the chemical names of the components of that essential oil or extract as referred to in Article 18(3)(b).

1.5. **Exemptions from labelling and packaging requirements**

1.5.1. ***Exemptions from Article 31 [(Article 29(1))]***

1.5.1.1. Where Article 29(1) applies, the label elements mentioned in Article 17 may be provided in one of the following ways:

(a) in fold-out labels; or

(b) on tie-on tags; or

(c) on an outer packaging.

1.5.1.2. The label on any inner packaging shall contain at least hazard pictograms, the product identifier referred to in Article 18 and name and telephone number of the supplier of the substance or mixture.

1.5.2. ***Exemptions from Article 17 [(Article 29(2)]***

1.5.2.1. *Labelling of packages where the contents do not exceed 125 ml*

1.5.2.1.1. The hazard statements and the precautionary statements linked to the hazard categories listed below may be omitted from the label elements required by Article 17 where:

(a) the contents of the package do not exceed 125 ml; and

(b) the substance or mixture is classified in one or more of the following hazard categories:

1) Oxidising gases of category 1;

2) Gases under pressure;

3) Flammable liquids of category 2 or 3;

4) Flammable solids of category 1 or 2;

5) Self-reactive substances or mixtures Types C to F;

6) Self-heating substances or mixtures of category 2;

7) Substances and mixtures which, in contact with water, emit flammable gases of categories 1, 2 or 3;

8) Oxidising liquids of category 2 or 3;

9) Oxidising solids of category 2 or 3;

10) Organic peroxides Types C to F;

11) Acute toxicity of category 4, if the substances or mixtures are not supplied to the general public;

12) Skin irritation of category 2;

13) Eye irritation of category 2;

14) Specific target organ toxicity — single exposure of category 2 or 3, if the substance or mixture is not supplied to the general public;

15) Specific target organ toxicity — repeated exposure of category 2, if the substance or mixture is not supplied to the general public;

16) Hazardous to the aquatic environment — Acute of category 1;

17) Hazardous to the aquatic environment — Chronic of category 1 or 2.

The exemptions for labelling of small packages of aerosols as flammable laid down in Directive 75/324/EEC shall apply to aerosol dispensers.

1.5.2.1.2. The precautionary statements linked to the hazard categories listed below may be omitted from the label elements required by Article 17 where:

(a) the contents of the package do not exceed 125 ml; and

(b) the substance or mixture is classified in one or more of the following hazard categories:

1) Flammable gases of category 2;

2) Reproductive toxicity: effects on or via lactation;

3) Hazardous to the aquatic environment — Chronic of category 3 or 4.

1.5.2.1.3. The pictogram, the hazard statement and the precautionary statement linked to the hazard categories listed below may be omitted from the label elements required by Article 17 where:

(a) the contents of the package do not exceed 125 ml; and

(b) the substance or mixture is classified in one or more of the following hazard categories:

1) Corrosive to metals.

1.5.2.2. *Labelling of soluble packaging for single use*

The label elements required by Article 17 may be omitted from soluble packaging intended for single use where:

(a) The content of each soluble packaging does not exceed a volume of 25 ml;

(b) The classification of the contents of the soluble packaging is exclusively one or more of the hazard categories in 1.5.2.1.1 (b); and

(c) The soluble packaging is contained within outer packaging that fully meets the requirements of Article 17.

1.5.2.3. Section 1.5.2.2 shall not apply to substances or mixtures within the scope of Directives 91/414/EEC or 98/8/EC.

2. PART 2: PHYSICAL HAZARDS

2.1. **Explosives**

2.1.1. ***Definitions***

2.1.1.1. The class of explosives comprises

(a) explosive substances and mixtures;

(b) explosive articles, except devices containing explosive substances or mixtures in such quantity or of such a character that their inadvertent or accidental ignition or initiation shall not cause any effect external to the device either by projection, fire, smoke, heat or loud noise; and

(c) substances, mixtures and articles not mentioned in points (a) and (b) which are manufactured with a view to producing a practical, explosive or pyrotechnic effect.

2.1.1.2. For the purposes of this Regulation the following definitions shall apply:

An explosive substance or mixture is a solid or liquid substance or mixture of substances which is in itself capable by chemical reaction of producing gas at such a temperature and pressure and at such a speed as to cause damage to the surroundings. Pyrotechnic substances are included even when they do not evolve gases.

A pyrotechnic substance or mixture is a substance or mixture of substances designed to produce an effect by heat, light, sound, gas or smoke or a combination of these as the result of non-detonative self-sustaining exothermic chemical reactions.

An unstable explosive is an explosive substance or mixture which is thermally unstable and/or too sensitive for normal handling, transport and use.

An explosive article is an article containing one or more explosive substances or mixtures.

A pyrotechnic article is an article containing one or more pyrotechnic substances or mixtures.

An intentional explosive is a substance, mixture or article which is manufactured with a view to producing a practical, explosive or pyrotechnic effect.

2.1.2. ***Classification criteria***

2.1.2.1. Substances, mixtures and articles of this class are classified as an unstable explosive on the basis of the flowchart in Figure 2.1.2. The test methods are described in Part I of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria.

2.1.2.2. Substances, mixtures and articles of this class, which are not classified as an unstable explosive, shall be assigned to one of the following six divisions depending on the type of hazard they present:

(a) Division 1.1 Substances, mixtures and articles which have a mass explosion hazard (a mass explosion is one which affects almost the entire quantity present virtually instantaneously);

(b) Division 1.2 Substances, mixtures and articles which have a projection hazard but not a mass explosion hazard;

(c) Division 1.3 Substances, mixtures and articles which have a fire hazard and either a minor blast hazard or a minor projection hazard or both, but not a mass explosion hazard:

(i) combustion of which gives rise to considerable radiant heat; or

(ii) which burn one after another, producing minor blast or projection effects or both;

(d) Division 1.4 Substances, mixtures and articles which present no significant hazard:

— substances, mixtures and articles which present only a small hazard in the event of ignition or initiation. The effects are largely confined to the package and no projection of fragments of appreciable size or range is to be expected. An external fire shall not cause virtually instantaneous explosion of almost the entire contents of the package;

(e) Division 1.5 Very insensitive substances or mixtures which have a mass explosion hazard:

— substances and mixtures which have a mass explosion hazard but are so insensitive that there is very little probability of initiation or of transition from burning to detonation under normal conditions;

(f) Division 1.6 Extremely insensitive articles which do not have a mass explosion hazard:

— articles which contain only extremely insensitive detonating substances or mixtures and which demonstrate a negligible probability of accidental initiation or propagation.

2.1.2.3. Explosives, which are not classified as an unstable explosive, shall be classified in one of the six divisions referred to in paragraph 2.1.2.2 of this Annex based on Test Series 2 to 8 in Part I of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria according to the results of the tests laid down in Table 2.1.1:

Table 2.1.1

Criteria for explosives

Category	Criteria
Unstable explosives or explosives of Divisions 1.1 to 1.6	For explosives of Divisions 1.1 to 1.6, the following are the core set of tests that need to be performed: Explosibility: according to UN Test Series 2 (section 12 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria). Intentional explosives ([1]) shall not be subject to UN Test Series 2. Sensitiveness: according to UN Test Series 3 (section 13 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria). Thermal stability: according to UN Test 3(c) (sub-section 13.6.1 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria). Further tests are necessary to allocate the correct Division.

([1]) This comprises substances, mixtures and articles which are manufactured with a view to producing a practical, explosive or pyrotechnic effect.

2.1.2.4. If explosives are unpackaged or repacked in packaging other than the original or similar packaging, they shall be retested.

2.1.3. ***Hazard Communication***

Label elements shall be used for substances, mixtures or articles meeting the criteria for classification in this hazard class in accordance with Table 2.1.2.

NOTE to Table 2.1.2: Unpackaged explosives or explosives repacked in packaging other than the original or similar packaging shall include all of the following label elements:

(a) the pictogram: exploding bomb;

(b) the signal word: 'Danger'; and

(c) the hazard statement: 'explosive; mass explosion hazard'

unless the hazard is shown to correspond to one of the hazard categories in Table 2.1.2, in which case the corresponding symbol, the signal word and/or the hazard statement shall be assigned.

Table 2.1.2

Label elements for explosives

Classification	Unstable Explosive	Division 1.1	Division 1.2	Division 1.3	Division 1.4	Division 1.5	Division 1.6
GHS Pictograms							
Signal Word	Danger	Danger	Danger	Danger	Warning	Danger	No signal word
Hazard Statement	H200: Unstable Explosive	H201: Explosive; mass explosion hazard	H202: Explosive; severe projection hazard	H203: Explosive; fire, blast or projection hazard	H204: Fire or projection hazard	H205: May mass explode in fire	No hazard statement
Precautionary Statement Prevention	P201 P202 P281	P210 P230 P240 P250 P280	P210 P230 P240 P250 P280	P210 P230 P240 P250 P280	P210 P240 P250 P280	P210 P230 P240 P250 P280	No precautionary statement
Precautionary Statement Response	P372 P373 P380	P370+P380 P372 P373	P370+P380 P372 P373	P370+P380 P372 P373	P370+P380 P372 P373	P370+P380 P372 P373	No precautionary statement
Precautionary Statement Storage	P401	P401	P401	P401	P401	P401	No precautionary statement
Precautionary Statement Disposal	P501	P501	P501	P501	P501	P501	No precautionary statement

2.1.4. ***Additional Classification Considerations***

2.1.4.1. The classification of substances, mixtures and articles in the explosives hazard class and further allocation to a division is a very complex, three step procedure. Reference to Part I of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria is necessary.

The first step is to ascertain whether the substance or mixture has explosive effects (Test Series 1). The second step is the acceptance procedure (Test Series 2 to 4) and the third step is the assignment to a hazard division (Test Series 5 to 7). The assessment whether a candidate for 'ammonium nitrate emulsion or suspension or gel, intermediate for blasting explosives (ANE)' is insensitive enough for inclusion as an oxidising liquid (section 2.13) or an oxidising solid (section 2.14) is answered by Test Series 8 tests.

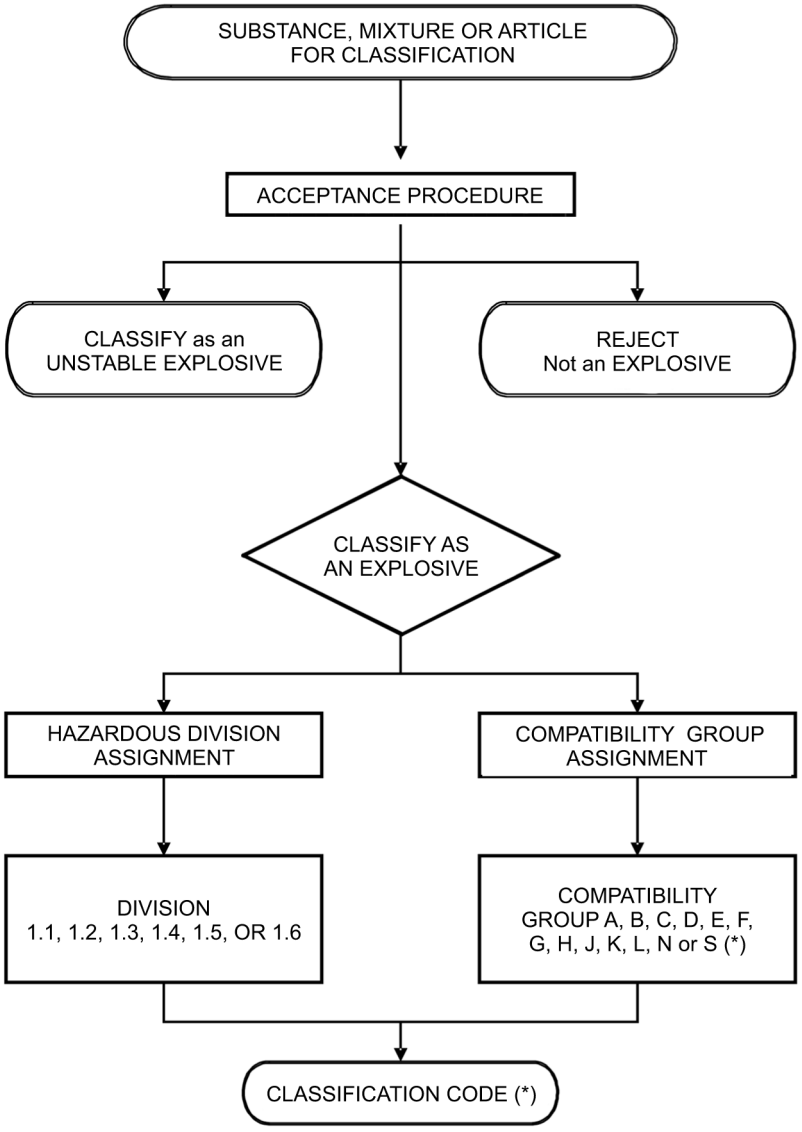
Explosive substances and mixtures wetted with water or alcohols, or diluted with other substances to suppress their explosive properties, may be treated differently in terms of classification and other hazard classes may apply, according to their physical properties (see also Annex II section 1.1.).

Certain physical hazards (due to explosive properties) are altered by dilution, as is the case for desensitised explosives, by inclusion in a mixture or article, packaging or other factors.

The classification procedure is set out in the following decision logic (see Figures 2.1.1 to 2.1.4).

Figure 2.1.1

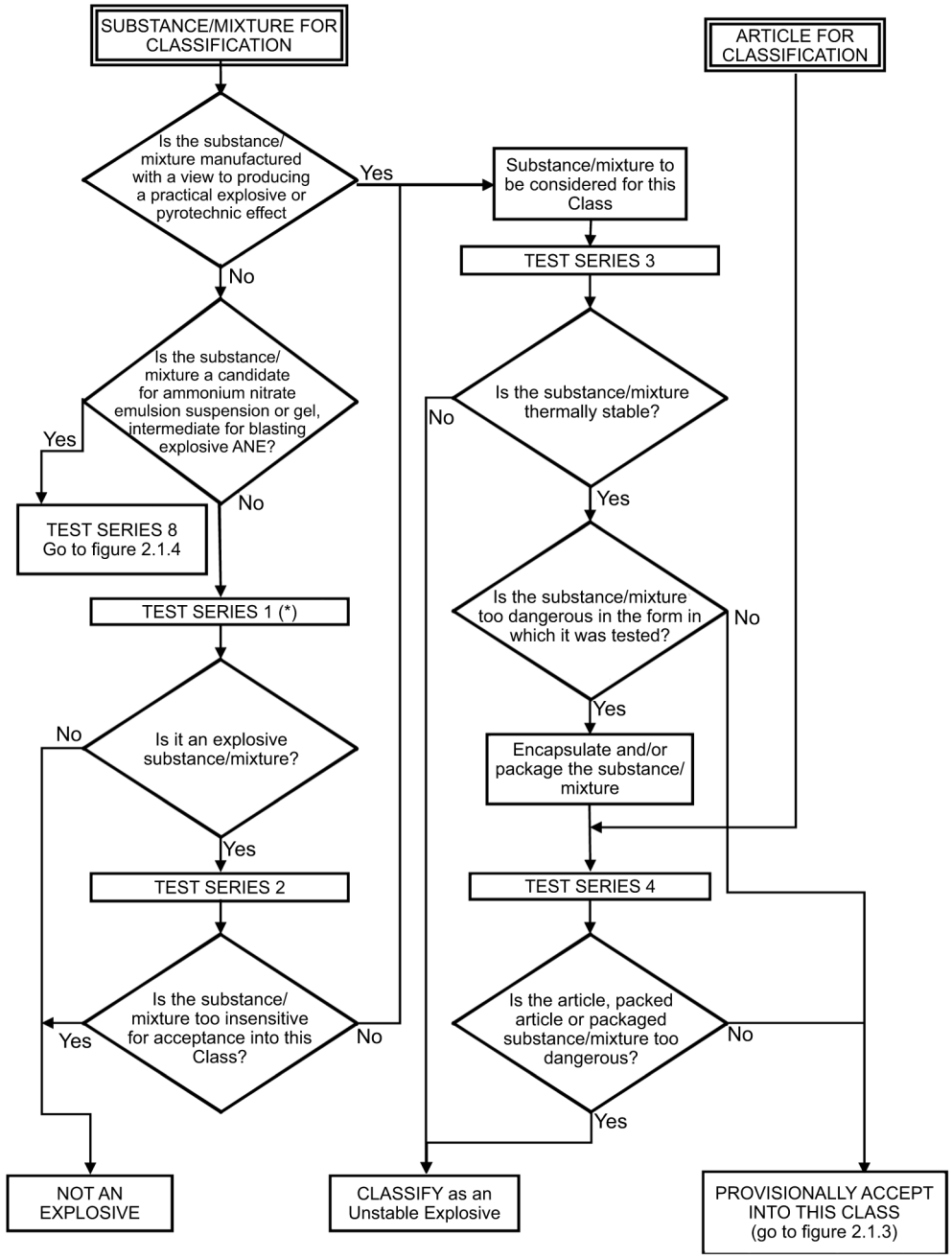
Overall scheme of the procedure for classifying a substance, mixture or article in the class of explosives (Class 1 for transport)



(*) see UN Recommendations on the Transport of Dangerous Goods, Model Regulations, 15th rev. ed, sub-section 2.1.2.

Figure 2.1.2

Procedure for provisional acceptance of a substance, mixture or article in the class of explosives (Class 1 for transport)



(*) For classification purposes, start with Test Series 2.

Figure 2.1.3

Procedure for assignment to a division in the class of explosives (Class 1 for transport)

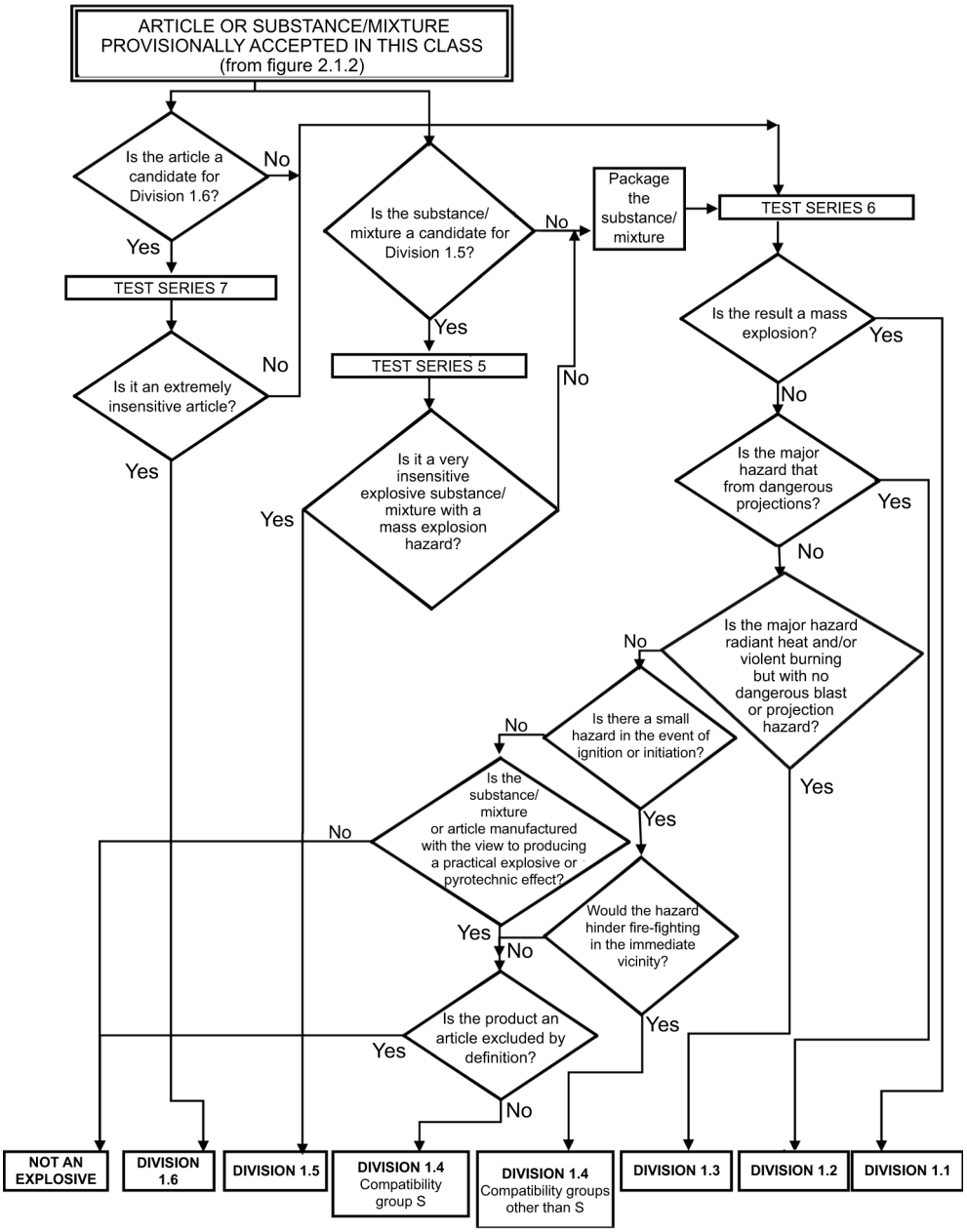
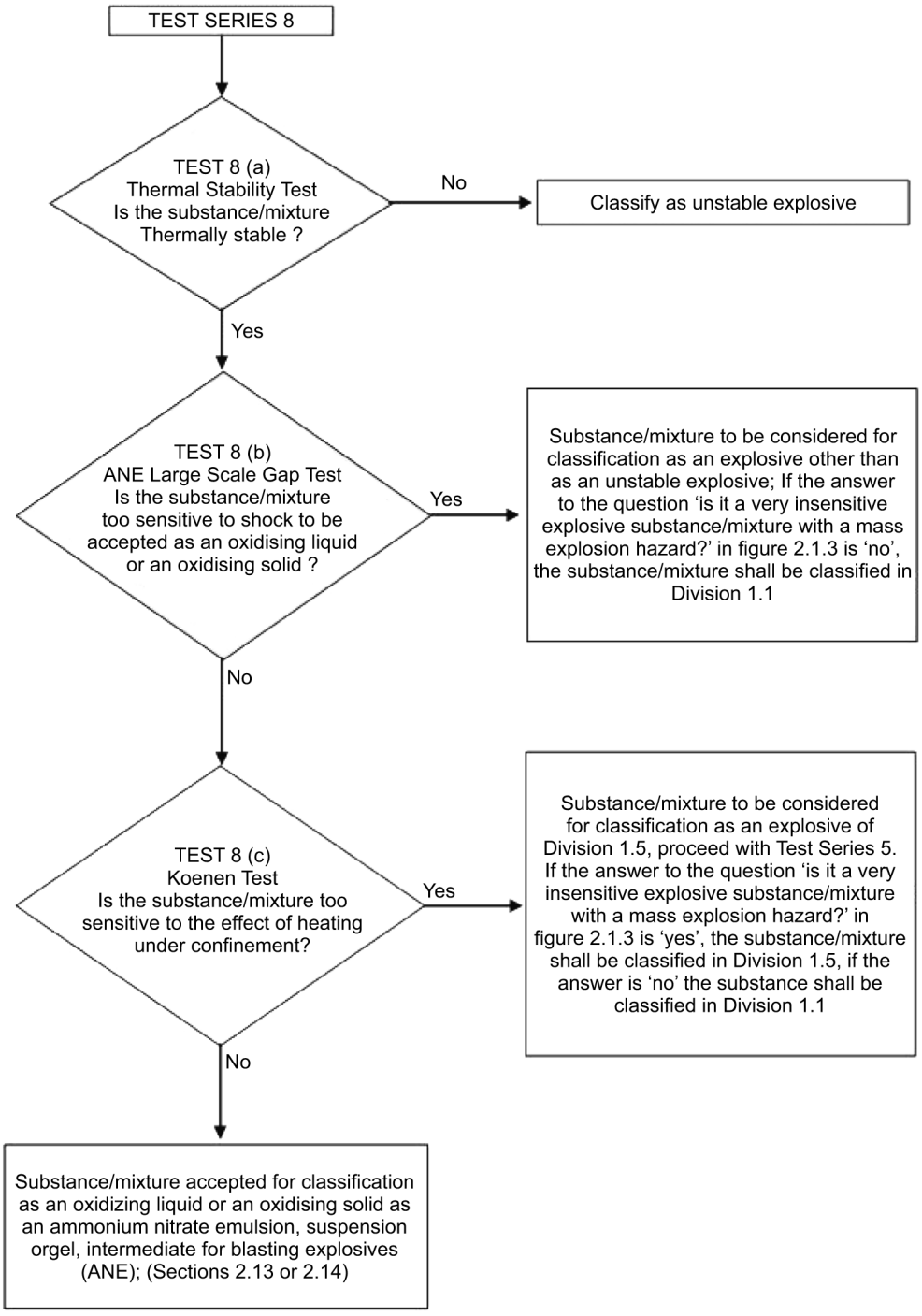


Figure 2.1.4

Procedure for classification of ammonium nitrate emulsions, suspensions or gels



2.1.4.2. *Screening procedure*

Explosive properties are associated with the presence of certain chemical groups in a molecule which can react to produce very rapid increases in temperature or pressure. The screening procedure is aimed at identifying the presence of such reactive groups and the potential for rapid energy release. If the screening procedure identifies the substance or mixture to be a potential explosive, the acceptance procedure (see section 10.3 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria) has to be performed.

Note

Neither a Series 1 type (a) propagation of detonation test nor a Series 2 type (a) test of sensitivity to detonative shock is required if the exothermic decomposition energy of organic materials is less than 800 J/g.

2.1.4.3. A substance or mixture shall not be classified as explosive if:

(a) There are no chemical groups associated with explosive properties present in the molecule. Examples of groups which may indicate explosive properties are given in Table A6.1 in Appendix 6 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria; or

(b) The substance contains chemical groups associated with explosive properties which include oxygen and the calculated oxygen balance is less than - 200;

The oxygen balance is calculated for the chemical reaction:

$C_xH_yO_z + [x + (y/4) - (z/2)]\ O_2 \rightarrow x\ CO_2 + (y/2)\ H_2O$

Using the formula:

Oxygen balance = - 1 600 [2x + (y/2)-z]/molecular weight;

(c) When the organic substance or a homogenous mixture of organic substances contains chemical groups associated with explosive properties but the exothermic decomposition energy is less than 500 J/g and the onset of exothermic decomposition is below 500 °C. The exothermic decomposition energy can be determined using a suitable calorimetric technique; or

(d) For mixtures of inorganic oxidising substances with organic material(s), the concentration of the inorganic oxidising substance is:

— less than 15 % by mass, if the oxidising substance is assigned to Categories 1 or 2;

— less than 30 % by mass, if the oxidising substance is assigned to Category 3.

2.1.4.4. In the case of mixtures containing any known explosives, the acceptance procedure has to be performed.

2.2. Flammable gases

2.2.1. *Definition*

Flammable gas means a gas or gas mixture having a flammable range with air at 20 °C and a standard pressure of 101,3 kPa.

2.2.2. *Classification criteria*

2.2.2.1. A flammable gas shall be classified in this class in accordance with Table 2.2.1:

Table 2.2.1

Criteria for flammable gases

Category	Criteria
1	Gases, which at 20 °C and a standard pressure of 101,3 kPa: (a) are ignitable when in a mixture of 13 % or less by volume in air; or (b) have a flammable range with air of at least 12 percentage points regardless of the lower flammable limit.
2	Gases, other than those of Category 1, which, at 20 °C and a standard pressure of 101,3 kPa, have a flammable range while mixed in air.

Note

For the classification of aerosols, see 2.3.

2.2.3. ***Hazard Communication***

Label elements shall be used for substances and mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.2.2.

Table 2.2.2

Label elements for flammable gases

Classification	Category 1	Category 2
GHS Pictogram		No pictogram
Signal Word	Danger	Warning
Hazard Statement	H220: Extremely flammable gas	H221: Flammable gas
Precautionary Statement Prevention	P210	P210
Precautionary Statement Response	P377 P381	P377 P381
Precautionary Statement Storage	P403	P403
Precautionary Statement Disposal		

2.2.4. ***Additional Classification Considerations***

2.2.4.1. Flammability shall be determined by tests or, for mixtures where there are sufficient data available, by calculation in accordance with the methods adopted by ISO (see ISO 10156 as amended, Gases and gas mixtures — Determination of fire potential and oxidising ability for the selection of cylinder valve outlet). Where insufficient data are available to use these methods, test method EN 1839 as amended (Determination of explosion limits of gases and vapours) can be used.

2.3. **Flammable aerosols**

2.3.1. ***Definitions***

Aerosols, this means aerosol dispensers, are any non-refillable receptacles made of metal, glass or plastics and containing a gas compressed, liquefied or dissolved under pressure, with or without a liquid, paste or powder, and fitted with a release device allowing the contents to be ejected as solid or liquid particles in suspension in a gas, as a foam, paste or powder or in a liquid state or in a gaseous state.

2.3.2. ***Classification criteria***

2.3.2.1. Aerosols shall be considered for classification as flammable in accordance with 2.3.2.2 if they contain any component which is classified as flammable according to the criteria contained in this Part, i.e.:

— Liquids with a flash point ≤ 93 °C, which includes Flammable Liquids according to section 2.6;

— Flammable gases (see 2.2);

— Flammable solids (see 2.7).

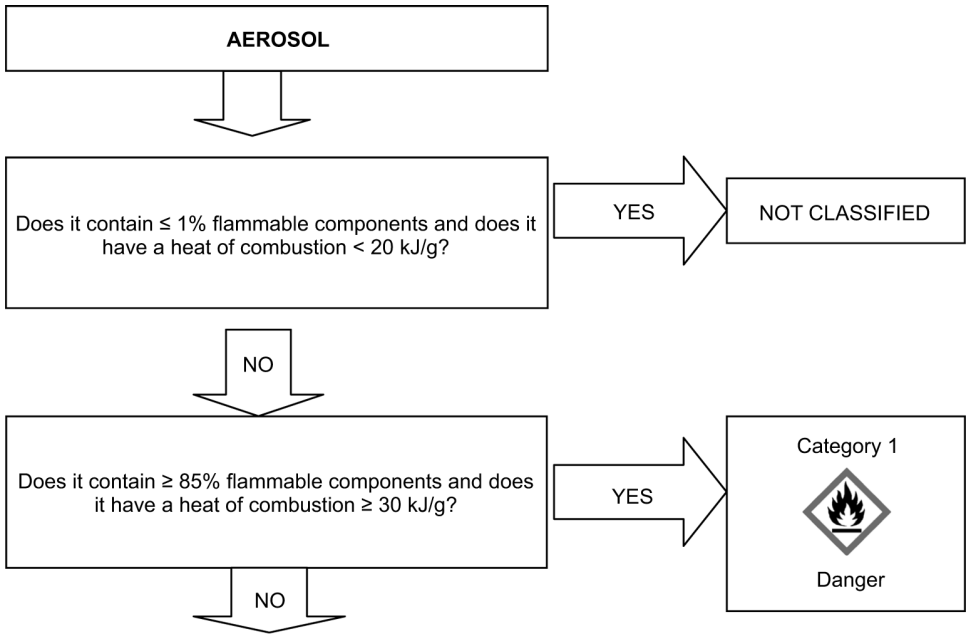
Note

Flammable components do not cover pyrophoric, self-heating or water-reactive substances and mixtures because such components are never used as aerosol contents.

2.3.2.2. A flammable aerosol shall be classified in one of the two categories for this Class on the basis of its components, of its chemical heat of combustion and, if applicable, of the results of the foam test (for foam aerosols) and of the ignition distance test and enclosed space test (for spray aerosols) in accordance with Figure 2.3.1 and the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, Part III, sub-sections 31.4, 31.5 and 31.6.

Figure 2.3.1

Figure 2.3.1(a) for flammable aerosols



For spray aerosols, go to decision logic 2.3.1 (b);

For foam aerosols, got to decision logic 2.3.1 (c).

Figure 2.3.1(b) for spray aerosols

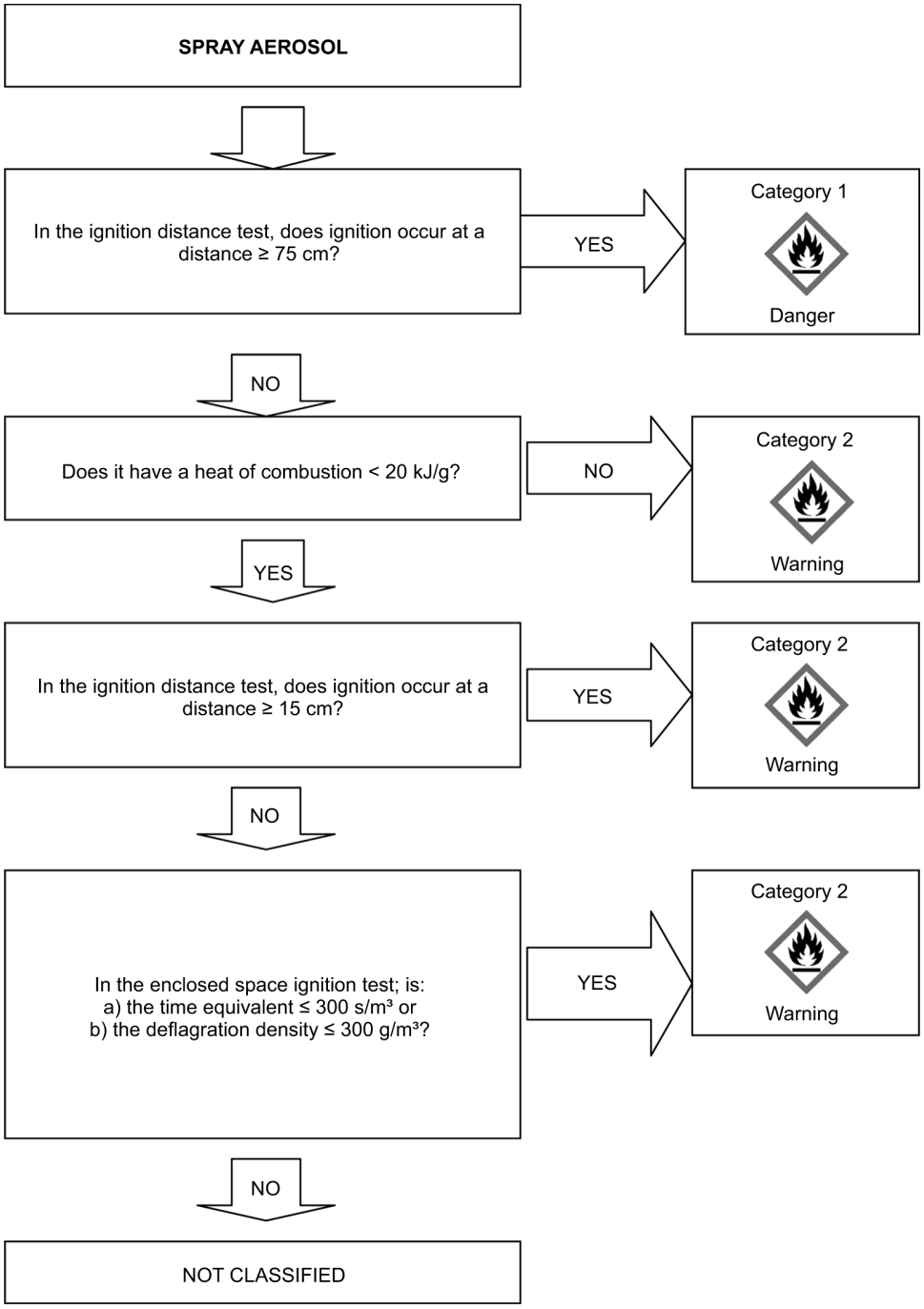
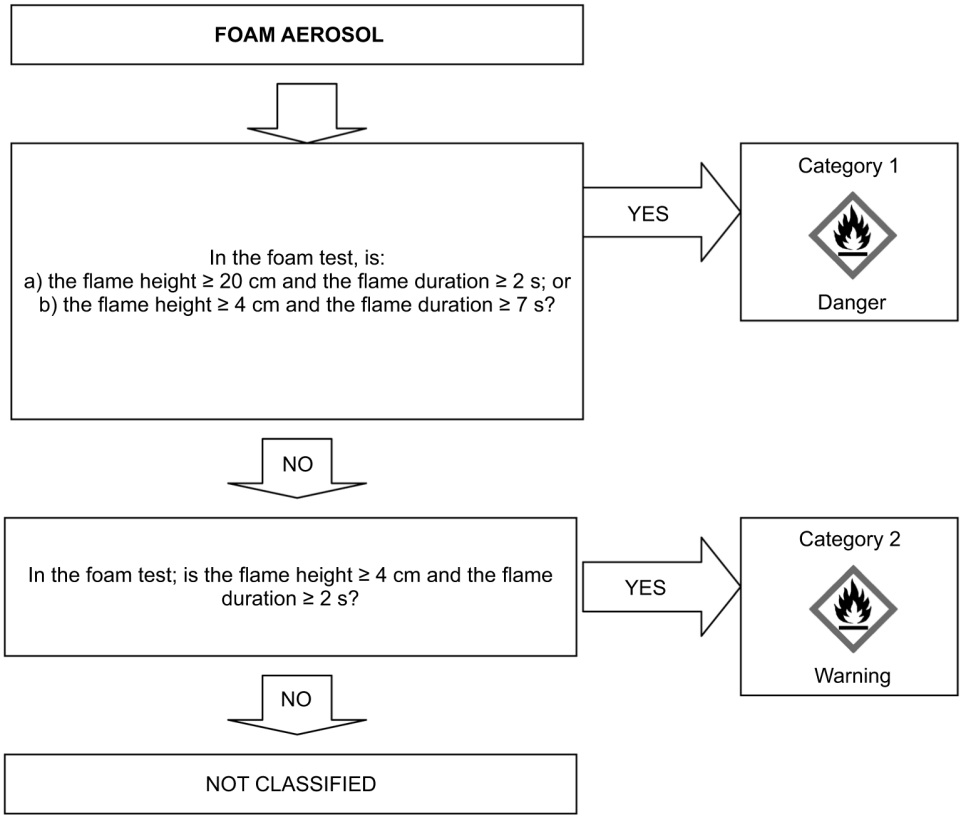


Figure 2.3.1(c) for foam aerosols



2.3.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.3.2.

Table 2.3.2

Label elements for flammable aerosols

Classification	Category 1	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H222: Extremely flammable aerosol	H223: Flammable aerosol
Precautionary Statement Prevention	P210 P211 P251	P210 P211 P251
Precautionary Statement Response		
Precautionary Statement Storage	P410 + P412	P410 + P412
Precautionary Statement Disposal		

2.3.4. ***Additional Classification Considerations***

2.3.4.1. The chemical heat of combustion (ΔH_c), in kilojoules per gram (kJ/g), is the product of the theoretical heat of combustion (ΔH_{comb}), and a combustion efficiency, usually less than 1,0 (a typical combustion efficiency is 0,95 or 95 %).

For a composite aerosol formulation, the chemical heat of combustion is the summation of the weighted heats of combustion for the individual components, as follows:

$$\Delta H_{c\ (product)} = \sum_{i}^{n} \left[w_i\ \% \times \Delta H_{c(i)} \right]$$

where:

ΔH_c = chemical heat of combustion (kJ/g);
w_i % = mass fraction of component i in the product;
$\Delta H_{c(i)}$ = specific heat of combustion (kJ/g)of component i in the product.

The chemical heats of combustion can be found in the literature, calculated or determined by tests (see ASTM D 240 as amended — Standard Test Methods for Heat of Combustion of Liquid Hydrocarbon Fuels by Bomb Calorimeter, EN/ISO 13943 as amended, 86.l to 86.3 — Fire safety — Vocabulary, and NFPA 30B as amended — Code for the Manufacture and Storage of Aerosol Products).

2.4. **Oxidising gases**

2.4.1. ***Definitions***

Oxidising gas means any gas or gas mixture which may, generally by providing oxygen, cause or contribute to the combustion of other material more than air does.

2.4.2. ***Classification criteria***

2.4.2.1. An oxidising gas shall be classified in a single category for this class in accordance with Table 2.4.1.:

Table 2.4.1

Criteria for oxidising gases

Category	Criteria
1	Any gas which may, generally by providing oxygen, cause or contribute to the combustion of other material more than air does.

Note

'Gases which cause or contribute to the combustion of other material more than air' means pure gases or gas mixtures with an oxidising power greater than 23,5 % as determined by a method specified in ISO 10156 as amended or 10156-2 as amended.

2.4.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.4.2.

Table 2.4.2

Label elements for oxidising gases

Classification	Category 1
GHS Pictogram	

Classification	Category 1
Signal Word	Danger
Hazard Statement	H270: May cause or intensify fire; oxidiser
Precautionary Statement Prevention	P220 P244
Precautionary Statement Response	P370 + P376
Precautionary Statement Storage	P403
Precautionary Statement Disposal	

2.4.4. ***Additional Classification Considerations***

To classify an oxidising gas, tests or calculation methods as described in ISO 10156 as amended, gases and gas mixtures — Determination of fire potential and oxidising ability for the selection of cylinder valve outlet and ISO 10156-2 as amended, gas cylinders — gases and gas mixtures — Determination of oxidising ability of toxic and corrosive gases and gas mixtures — shall be performed.

2.5. **Gases under pressure**

2.5.1. ***Definition***

2.5.1.1. Gases under pressure are gases which are contained in a receptacle at a pressure of 200 kPa (gauge) or more, or which are liquefied or liquefied and refrigerated.

They comprise compressed gases, liquefied gases, dissolved gases and refrigerated liquefied gases.

2.5.1.2. The critical temperature is the temperature above which a pure gas cannot be liquefied, regardless of the degree of compression.

2.5.2. ***Classification criteria***

Gases shall be classified, according to their physical state when packaged, in one of four groups in accordance with Table 2.5.1:

Table 2.5.1

Criteria for gases under pressure

Group	Criteria
Compressed gas	A gas which when packaged under pressure is entirely gaseous at - 50 °C; including all gases with a critical temperature ≤ - 50 °C.
Liquefied gas	A gas which, when packaged under pressure, is partially liquid at temperatures above - 50 °C. A distinction is made between: (i) high pressure liquefied gas: a gas with a critical temperature between - 50 °C and + 65 °C; and (ii) low pressure liquefied gas: a gas with a critical temperature above + 65 °C.
Refrigerated liquefied gas	A gas which when packaged is made partially liquid because of its low temperature.
Dissolved gas	A gas which when packaged under pressure is dissolved in a liquid phase solvent.

2.5.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.5.2.

Table 2.5.2

Label elements for gases under pressure

Classification	Compressed gas	Liquefied gas	Refrigerated liquefied gas	Dissolved gas
GHS Pictograms				
Signal Word	Warning	Warning	Warning	Warning
Hazard Statement	H280: Contains gas under pressure; may explode if heated	H280: Contains gas under pressure; may explode if heated	H281: Contains refrigerated gas; may cause cryogenic burns or injury	H280: Contains gas under pressure; may explode if heated
Precautionary Statement Prevention			P282	
Precautionary Statement Response			P336 P315	
Precautionary Statement Storage	P410 + P403	P410 + P403	P403	P410 + P403
Precautionary Statement Disposal				

2.5.4. ***Additional Classification Considerations***

For this group of gases, the following information is required to be known:

— the vapour pressure at 50 °C;

— the physical state at 20 °C at standard ambient pressure;

— the critical temperature.

Data can be found in the literature, calculated or determined by testing. Most pure gases are already classified in the UN Recommendations on the Transport of Dangerous Goods, Model Regulations.

2.6. **Flammable liquids**

2.6.1. ***Definition***

Flammable liquid means a liquid having a flash point of not more than 60 °C.

2.6.2. ***Classification criteria***

2.6.2.1. A flammable liquid shall be classified in one of the three categories for this class in accordance with Table 2.6.1:

Table 2.6.1

Criteria for flammable liquids

Category	Criteria
1	Flash point < 23 °C and initial boiling point ≤ 35 °C
2	Flash point < 23 °C and initial boiling point > 35 °C
3	Flash point ≥ 23 °C and ≤ 60 °C ([1])

([1]) For the purpose of this Regulation gas oils, diesel and light heating oils having a flash point between ≥ 55 °C and ≤ 75 °C may be regarded as Category 3.

2.6.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.6.2.

Table 2.6.2

Label elements for flammable liquids

Classification	Category 1	Category 2	Category 3
GHS Pictograms			
Signal Word	Danger	Danger	Warning
Hazard Statement	H224: Extremely flammable liquid and vapour	H225: Highly flammable liquid and vapour	H226: Flammable liquid and vapour
Precautionary Statement Prevention	P210 P233 P240 P241 P242 P243 P280	P210 P233 P240 P241 P242 P243 P280	P210 P233 P240 P241 P242 P243 P280
Precautionary Statement Response	P303 + P361 + P353 P370 + P378	P303 + P361 + P353 P370 + P378	P303 + P361 + P353 P370 + P378
Precautionary Statement Storage	P403 + P235	P403 + P235	P403 + P235
Precautionary Statement Disposal	P501	P501	P501

2.6.4. ***Additional Classification Considerations***

2.6.4.1. For the classification of flammable liquids data on flash point and initial boiling point are needed. Data can be determined by testing, found in literature or calculated. If data are not available, the flash point and the initial boiling point shall be determined through testing. For flash point determination a closed-cup method shall be used.

2.6.4.2. In the case of mixtures ([1]) containing known flammable liquids in defined concentrations, although they may contain non-volatile components e.g. polymers, additives, the flash point need not be determined experimentally if the calculated flash point of the mixture, using the method given in 2.6.4.3, is at least 5 °C ([2]) greater than the relevant classification criterion and provided that:

(a) the composition of the mixture is accurately known (if the material has a specified range of composition, the composition with the lowest calculated flash point shall be selected for assessment);

(b) the lower explosion limit of each component is known (an appropriate correlation has to be applied when these data are extrapolated to other temperatures than test conditions) as well as a method for calculating the lower explosion limit;

(c) the temperature dependence of the saturated vapour pressure and of the activity coefficient is known for each component as present in the mixture;

(d) the liquid phase is homogeneous.

2.6.4.3. One suitable method is described in Gmehling and Rasmussen (Ind. Eng. Fundament, 21, 186, (1982)). For a mixture containing non-volatile components the flash point is calculated from the volatile components. It is considered that a non-volatile component only slightly decreases the partial pressure of the solvents and the calculated flash point is only slightly below the measured value.

2.6.4.4. Possible test methods for determining the flash point of flammable liquids are listed in Table 2.6.3.

Table 2.6.3

Methods for determining the flash point of flammable liquids

European standards:	EN ISO 1516 as amended Determination of flash/no flash — Closed cup equilibrium method
	EN ISO 1523 as amended Determination of flash point — Closed cup equilibrium method
	EN ISO 2719 as amended Determination of flash point — Pensky-Martens closed cup method
	EN ISO 3679 as amended Determination of flash point — Rapid equilibrium closed cup method
	EN ISO 3680 as amended Determination of flash/no flash — Rapid equilibrium closed cup method
	EN ISO 13736 as amended Petroleum products and other liquids — Determination of flash point — Abel closed cup method
National standards:	
Association française de normalisation, AFNOR:	NF M07-036 as amended Détermination du point d'éclair — Vase clos Abel-Pensky (identical to DIN 51755)
British Standards Institute,	BS 2000 Part 170 as amended (identical to EN ISO 13736)
Deutsches Institut für Normung	DIN 51755 (flash points below 65 C) as amended Prüfung von Mineralölen und anderen brennbaren Flüssigkeiten; Bestimmung des Flammpunktes im geschlossenen Tiegel, nach Abel-Pensky (identical to NF M07-036)

([1]) To date, the calculation method has been validated for mixtures containing up to 6 volatile components. These components may be flammable liquids like hydrocarbons, ethers, alcohols, esters (except acrylates), and water. It is however not yet validated for mixtures containing halogenated sulphurous, and/or phosphoric compounds as well as reactive acrylates.

([2]) If the calculated flash point is less than 5 °C greater than the relevant classification criterion, the calculation method may not be used and the flash point should be determined experimentally.

2.6.4.5. Liquids with a flash point of more than 35 °C need not be classified in Category 3 if negative results have been obtained in the sustained combustibility test L.2, Part III, section 32 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria.

2.7. **Flammable solids**

2.7.1. ***Definition***

2.7.1.1. A flammable solid means a solid which is readily combustible, or may cause or contribute to fire through friction.

Readily combustible solids are powdered, granular, or pasty substances or mixtures which are dangerous if they can be easily ignited by brief contact with an ignition source, such as a burning match, and if the flame spreads rapidly.

2.7.2. ***Classification criteria***

2.7.2.1. Powdered, granular or pasty substances or mixtures (except powders of metals or metal alloys — see 2.7.2.2) shall be classified as readily combustible solids when the time of burning of one or more of the test runs, performed in accordance with the test method described in Part III, sub-section 33.2.1, of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, is less than 45 seconds or the rate of burning is more than 2,2 mm/s.

2.7.2.2. Powders of metals or metal alloys shall be classified as flammable solids when they can be ignited and the reaction spreads over the whole length of the sample in 10 minutes or less.

2.7.2.3. A flammable solid shall be classified in one of the two categories for this class using Method N.1 as described in 33.2.1 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria in accordance with Table 2.7.1:

Table 2.7.1

Criteria for flammable solids

Category	Criteria
1	Burning rate test Substances and mixtures other than metal powders: (a) wetted zone does not stop fire and (b) burning time < 45 seconds or burning rate > 2,2 mm/s Metal powders burning time ≤ 5 minutes
2	Burning rate test Substances and mixtures other than metal powders: (a) wetted zone stops the fire for at least 4 minutes and (b) burning time < 45 seconds or burning rate > 2,2 mm/s Metal powders burning time > 5 minutes and ≤ 10 minutes

Note

The test shall be performed on the substance or mixture in its physical form as presented. If, for example, for the purposes of supply or transport, the same chemical is to be presented in a physical form different from that which was tested and which is considered likely to materially alter its performance in a classification test, the substance shall also be tested in the new form.

2.7.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.7.2.

Table 2.7.2

Label elements for flammable solids

Classification	Category 1	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H228: Flammable Solid	H228: Flammable Solid
Precautionary Statement Prevention	P210 P240 P241 P280	P210 P240 P241 P280
Precautionary Statement Response	P370 + P378	P370 + P378
Precautionary Statement Storage		
Precautionary Statement Disposal		

2.8. **Self-reactive substances and mixtures**

2.8.1. ***Definition***

2.8.1.1. Self-reactive substances or mixtures are thermally unstable liquid or solid substances or mixtures liable to undergo a strongly exothermic decomposition even without participation of oxygen (air). This definition excludes substances and mixtures classified according to this Part as explosives, organic peroxides or as oxidising.

2.8.1.2. A self-reactive substance or mixture is regarded as possessing explosive properties when in laboratory testing the formulation is liable to detonate, to deflagrate rapidly or to show a violent effect when heated under confinement.

2.8.2. ***Classification criteria***

2.8.2.1. Any self-reactive substance or mixture shall be considered for classification in this class as a self-reactive substance or mixture unless:

(a) they are explosives, according to the criteria given in 2.1;

(b) they are oxidising liquids or solids, according to the criteria given in 2.13 or 2.14, except that mixtures of oxidising substances, which contain 5 % or more of combustible organic substances shall be classified as self-reactive substances according to the procedure defined in 2.8.2.2;

(c) they are organic peroxides, according to the criteria given in 2.15;

(d) their heat of decomposition is less than 300 J/g; or

(e) their self-accelerating decomposition temperature (SADT) is greater than 75 °C for a 50 kg package [1].

[1] See UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, sub-sections 28.1, 28.2, 28.3 and Table 28.3.

2.8.2.2. Mixtures of oxidising substances, meeting the criteria for classification as oxidising substances, which contain 5 % or more of combustible organic substances and which do not meet the criteria mentioned in (a), (c), (d) or (e) in 2.8.2.1, shall be subjected to the self-reactive substances classification procedure;

Such a mixture showing the properties of a self-reactive substance type B to F (see 2.8.2.3) shall be classified as a self-reactive substance.

Where the test is conducted in the package form and the packaging is changed, a further test shall be conducted where it is considered that the change in packaging will affect the outcome of the test.

2.8.2.3. Self-reactive substances and mixtures shall be classified in one of the seven categories of 'types A to G' for this class, according to the following principles:

(a) any self-reactive substance or mixture which can detonate or deflagrate rapidly, as packaged, shall be defined as self-reactive substance TYPE A;

(b) any self-reactive substance or mixture possessing explosive properties and which, as packaged, neither detonates nor deflagrates rapidly, but is liable to undergo a thermal explosion in that package shall be defined as self-reactive substance TYPE B;

(c) any self-reactive substance or mixture possessing explosive properties when the substance or mixture as packaged cannot detonate or deflagrate rapidly or undergo a thermal explosion shall be defined as self-reactive substance TYPE C;

(d) any self-reactive substance or mixture which in laboratory testing:

(i) detonates partially, does not deflagrate rapidly and shows no violent effect when heated under confinement; or

(ii) does not detonate at all, deflagrates slowly and shows no violent effect when heated under confinement; or

(iii) does not detonate or deflagrate at all and shows a medium effect when heated under confinement;

shall be defined as self-reactive substance TYPE D;

(e) any self-reactive substance or mixture which, in laboratory testing, neither detonates nor deflagrates at all and shows low or no effect when heated under confinement shall be defined as self-reactive substance TYPE E;

(f) any self-reactive substance or mixture which, in laboratory testing, neither detonates in the cavitated state nor deflagrates at all and shows only a low or no effect when heated under confinement as well as low or no explosive power shall be defined as self-reactive substance TYPE F;

(g) any self-reactive substance or mixture which, in laboratory testing, neither detonates in the cavitated state nor deflagrates at all and shows no effect when heated under confinement nor any explosive power, provided that it is thermally stable (SADT is 60 °C to 75 °C for a 50 kg package), and, for liquid mixtures, a diluent having a boiling point not less than 150 °C is used for desensitisation shall be defined as self-reactive substance TYPE G. If the mixture is not thermally stable or a diluent having a boiling point less than 150 °C is used for desensitisation, the mixture shall be defined as self-reactive substance TYPE F.

Where the test is conducted in the package form and the packaging is changed, a further test shall be conducted where it is considered that the change in packaging will affect the outcome of the test.

2.8.2.4. *Criteria for temperature control*

Self-reactive substances need to be subjected to temperature control if their SADT is less than or equal to 55 °C. Test methods for determining the SADT as well as the derivation of control and emergency temperatures are given in, Part II, section 28 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria. The test selected shall be conducted in a manner which is representative, both in size and material, of the package.

2.8.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.8.1.

Table 2.8.1

Label elements for self-reactive substances and mixtures

Classification	Type A	Type B	Type C & D	Type E & F	Type G
GHS Pictograms					There are no label elements allocated to this hazard category
Signal Word	Danger	Danger	Danger	Warning	
Hazard Statement	H240: Heating may cause an explosion	H241: Heating may cause a fire or explosion	H242: Heating may cause a fire	H242: Heating may cause a fire	
Precautionary Statement Prevention	P210 P220 P234 P280	P210 P220 P234 P280	P210 P220 P234 P280	P210 P220 P234 P280	
Precautionary Statement Response	P370 + P378 P370 + P380 + P375	P370 + P378 P370 + P380 + P375	P370 + P378	P370 + P378	
Precautionary Statement Storage	P403 + P235 P411 P420	P403 + P235 P411 P420	P403 + P235 P411 P420	P403 + P235 P411 P420	
Precautionary Statement Disposal	P501	P501	P501	P501	

Type G has no hazard communication elements assigned but shall be considered for properties belonging to other hazard classes.

2.8.4. ***Additional Classification Considerations***

2.8.4.1. The properties of self-reactive substances or mixtures which are decisive for their classification shall be determined experimentally. The classification of a self reactive substance or mixture shall be performed in accordance with test series A to H as described in Part II of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria. The procedure for classification is described in Figure 2.8.1.

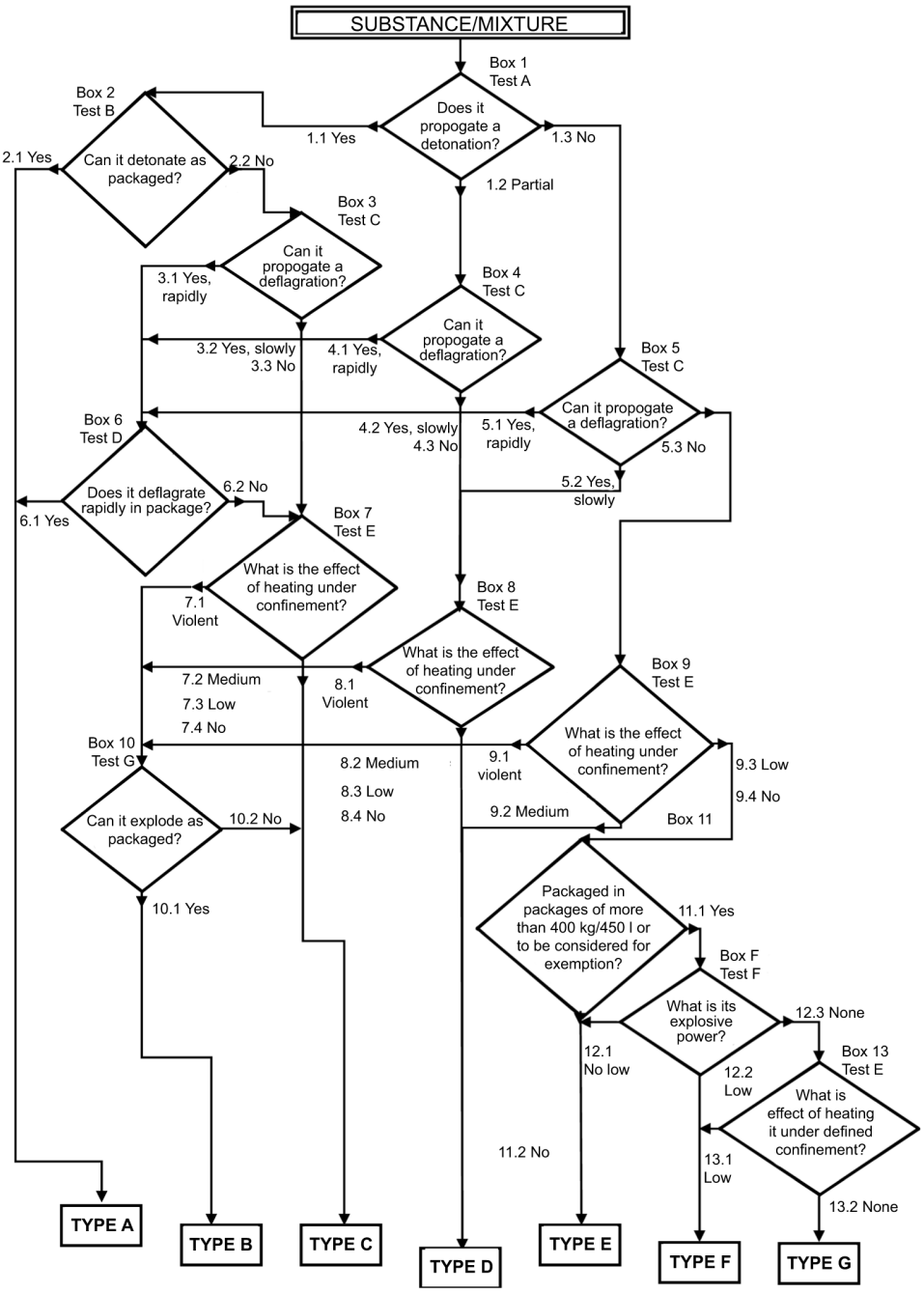
2.8.4.2. The classification procedures for self-reactive substances and mixtures need not be applied if:

(a) There are no chemical groups present in the molecule associated with explosive or self reactive properties. Examples of such groups are given in Tables A6.1 and A6.2 in Appendix 6 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria; or

(b) For a single organic substance or a homogeneous mixture of organic substances, the estimated SADT for a 50 kg package is greater than 75 °C or the exothermic decomposition energy is less than 300J/g. The onset temperature and decomposition energy can be estimated using a suitable calorimetric technique (see Part II, sub-section 20.3.3.3 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria).

Figure 2.8.1

Self-reactive substances and mixtures



2.9. **Pyrophoric liquids**

2.9.1. ***Definition***

Pyrophoric liquid means a liquid substance or mixture which, even in small quantities, is liable to ignite within five minutes after coming into contact with air.

2.9.2. ***Classification criteria***

2.9.2.1. A pyrophoric liquid shall be classified in a single category for this class using test N.3 in Part III, sub-section 33.3.1.5 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria according to Table 2.9.1:

Table 2.9.1

Criteria for pyrophoric liquids

Category	Criteria
1	The liquid ignites within 5 min when added to an inert carrier and exposed to air, or it ignites or chars a filter paper on contact with air within 5 min.

2.9.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.9.2.

Table 2.9.2

Label elements for pyrophoric liquids

Classification	Category 1
GHS Pictogram	
Signal Word	Danger
Hazard Statement	H250: Catches fire spontaneously if exposed to air
Precautionary Statement Prevention	P210 P222 P280
Precautionary Statement Response	P302 + P334 P370 + P378
Precautionary Statement Storage	P422
Precautionary Statement Disposal	

2.9.4. ***Additional Classification Considerations***

2.9.4.1. The classification procedure for pyrophoric liquids need not be applied when experience in manufacture or handling shows that the substance or mixture does not ignite spontaneously on coming into contact with air at normal temperatures (i.e. the substance is known to be stable at room temperature for prolonged periods of time (days)).

2.10. **Pyrophoric solids**

2.10.1. ***Definition***

Pyrophoric solid means a solid substance or mixture which, even in small quantities, is liable to ignite within five minutes after coming into contact with air.

2.10.2. ***Classification criteria***

2.10.2.1. A pyrophoric solid shall be classified in a single category for this class using test N.2 in Part III, sub-section 33.3.1.4 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria in accordance with Table 2.10.1:

Table 2.10.1

Criteria for pyrophoric solids

Category	Criteria
1	The solid ignites within 5 minutes of coming into contact with air.

Note

The test shall be performed on the substance or mixture in its physical form as presented. If, for example, for the purposes of supply or transport, the same chemical is to be presented in a physical form different from that which was tested and which is considered likely to materially alter its performance in a classification test, the substance shall also be tested in the new form.

2.10.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.10.2.

Table 2.10.2

Label elements for pyrophoric solids

Classification	Category 1
GHS Pictogram	
Signal Word	Danger
Hazard Statement	H250: Catches fire spontaneously if exposed to air
Precautionary Statement Prevention	P210 P222 P280
Precautionary Statement Response	P335 + P334 P370 +P378
Precautionary Statement Storage	P422
Precautionary Statement Disposal	

2.10.4. ***Additional Classification Considerations***

2.10.4.1. The classification procedure for pyrophoric solids need not be applied when experience in manufacture or handling shows that the substance or mixture does not ignite spontaneously on coming into contact with air at normal temperatures (i.e. the substance is known to be stable at room temperature for prolonged periods of time (days)).

2.11. **Self-heating substances and mixtures**

2.11.1. ***Definition***

2.11.1.1. A self-heating substance or mixture is a liquid or solid substance or mixture, other than a pyrophoric liquid or solid, which, by reaction with air and without energy supply, is liable to self-heat; this substance or mixture differs from a pyrophoric liquid or solid in that it will ignite only when in large amounts (kilograms) and after long periods of time (hours or days).

2.11.1.2. Self-heating of substances or mixtures, leading to spontaneous combustion, is caused by reaction of the substance or mixture with oxygen (in the air) and the heat developed not being conducted away rapidly enough to the surroundings. Spontaneous combustion occurs when the rate of heat production exceeds the rate of heat loss and the auto-ignition temperature is reached.

2.11.2. ***Classification criteria***

2.11.2.1. A substance or mixture shall be classified as a self-heating substance or mixture of this class, if in the tests performed in accordance with the test method given in the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, Part III, sub-section 33.3.1.6:

(a) a positive result is obtained using a 25 mm cube sample at 140 °C;

(b) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a negative result is obtained in a test using a 100 mm cube sample at 120 °C and the substance or mixture is to be packed in packages with a volume of more than 3 m^3;

(c) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a negative result is obtained in a test using a 100 mm cube sample at 100 °C and the substance or mixture is to be packed in packages with a volume of more than 450 litres;

(d) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a positive result is obtained in a test using a 100 mm cube sample at 100 °C.

2.11.2.2. A self-heating substance or mixture shall be classified in one of the two categories for this class if, in a test performed in accordance with test method N.4 in Part III, sub-section 33.3.1.6 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, the result meets the criteria according to Table 2.11.1:

Table 2.11.1

Criteria for self-heating substances and mixtures

Category	Criteria
1	A positive result is obtained in a test using a 25 mm sample cube at 140 °C
2	(a) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a negative result is obtained in a test using a 25 mm cube sample at 140 °C and the substance or mixture is to be packed in packages with a volume of more than 3 m^3; or (b) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a negative result is obtained in a test using a 25 mm cube sample at 140 °C, a positive result is obtained in a test using a 100 mm cube sample at 120 °C and the substance or mixture is to be packed in packages with a volume of more than 450 litres; or (c) a positive result is obtained in a test using a 100 mm sample cube at 140 °C and a negative result is obtained in a test using a 25 mm cube sample at 140 °C and a positive result is obtained in a test using a 100 mm cube sample at 100 °C.

Note

The test shall be performed on the substance or mixture in its physical form as presented. If, for example, for the purposes of supply or transport, the same chemical is to be presented in a physical form different from that which was tested and which is considered likely to materially alter its performance in a classification test, the substance shall also be tested in the new form.

2.11.2.3. Substances and mixtures with a temperature of spontaneous combustion higher than 50 °C for a volume of 27 m^3 shall not be classified as a self-heating substance or mixture.

2.11.2.4. Substances and mixtures with a spontaneous ignition temperature higher than 50 °C for a volume of 450 litres shall not be assigned to Category 1 of this class.

2.11.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.11.2.

Table 2.11.2

Label elements for self-heating substances and mixtures

Classification	Category 1	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H251: Self-heating; may catch fire	H252: Self-heating in large quantities; may catch fire
Precautionary Statement Prevention	P235 + P410 P280	P235 + P410 P280
Precautionary Statement Response		
Precautionary Statement Storage	P407 P413 P420	P407 P413 P420
Precautionary Statement Disposal		

2.11.4. ***Additional Classification Considerations***

2.11.4.1. For detailed schemes for the decision logic for classification and the tests to be carried out for ascertaining the different categories, see Figure 2.11.1.

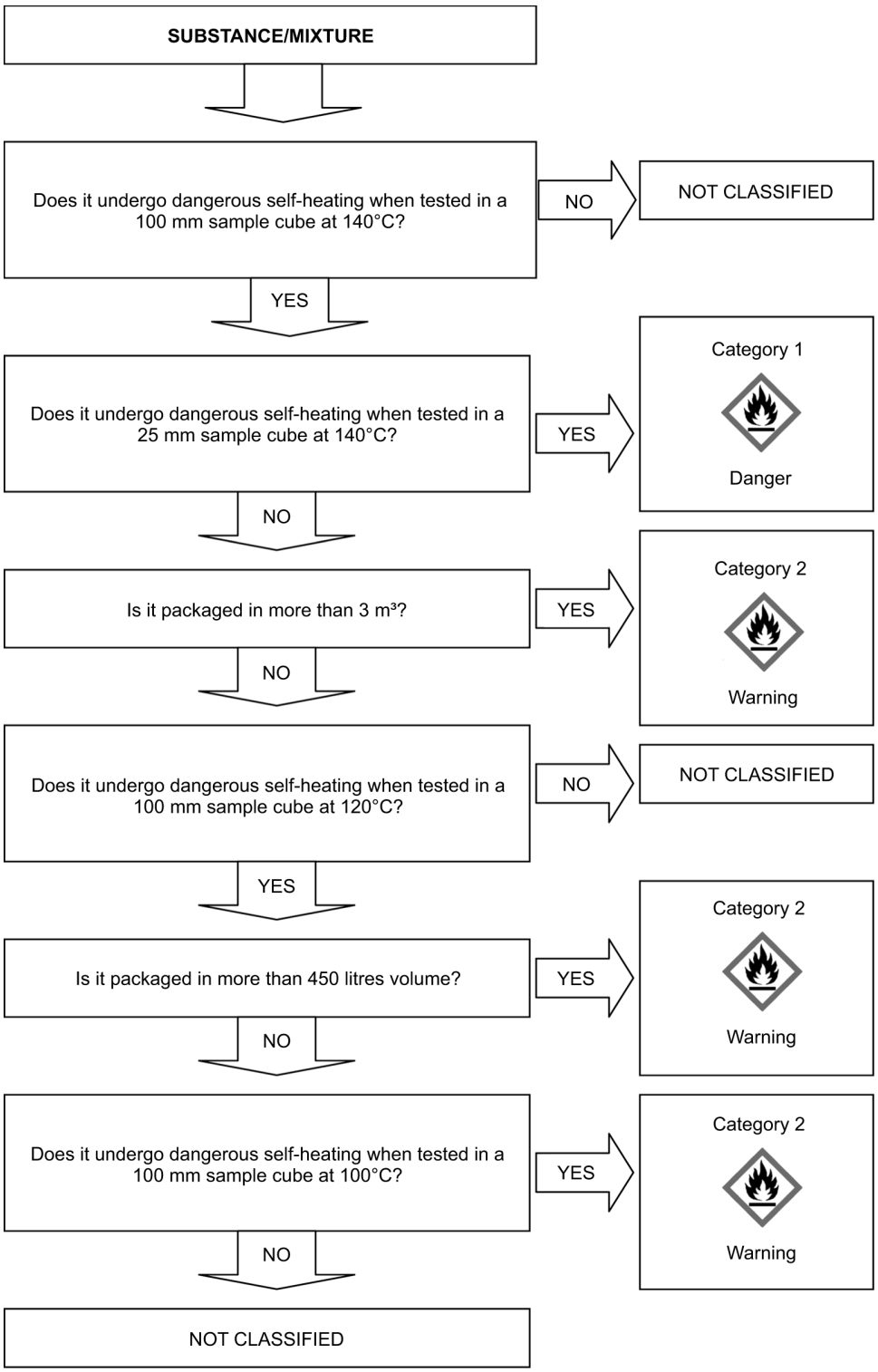
2.11.4.2. The classification procedure for self-heating substances or mixtures need not be applied if the results of a screening test can be adequately correlated with the classification test and an appropriate safety margin is applied. Examples of screening tests are:

(a) The Grewer Oven test (VDI guideline 2263, Part 1, 1990, Test methods for the Determination of the Safety Characteristics of Dusts) with an onset temperature 80 K above the reference temperature for a volume of 1 l;

(b) The Bulk Powder Screening Test (Gibson, N. Harper, D.J. Rogers, R.Evaluation of the fire and explosion risks in drying powders, Plant Operations Progress, 4 (3), 181-189, 1985) with an onset temperature 60 K above the reference temperature for a volume of 1 l.

Figure 2.11.1.

Self-heating substances and mixtures



2.12. **Substances and mixtures which in contact with water emit flammable gases**

2.12.1. ***Definition***

Substances or mixtures which, in contact with water, emit flammable gases means solid or liquid substances or mixtures which, by interaction with water, are liable to become spontaneously flammable or to give off flammable gases in dangerous quantities.

2.12.2. ***Classification criteria***

2.12.2.1. A substance or mixture which, in contact with water, emits flammable gases shall be classified in one of the three categories for this class, using test N.5 in Part III, sub-section 33.4.1.4 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, in accordance with Table 2.12.1:

Table 2.12.1

Criteria for substances or mixtures which in contact with water emit flammable gases

Category	Criteria
1	Any substance or mixture which reacts vigorously with water at ambient temperatures and demonstrates generally a tendency for the gas produced to ignite spontaneously, or which reacts readily with water at ambient temperatures such that the rate of evolution of flammable gas is equal to or greater than 10 litres per kilogram of substance over any one minute.
2	Any substance or mixture which reacts readily with water at ambient temperatures such that the maximum rate of evolution of flammable gas is equal to or greater than 20 litres per kilogram of substance per hour, and which does not meet the criteria for Category 1.
3	Any substance or mixture which reacts slowly with water at ambient temperatures such that the maximum rate of evolution of flammable gas is equal to or greater than 1 litre per kilogram of substance per hour, and which does not meet the criteria for Categories 1 and 2.

Note:

The test shall be performed on the substance or mixture in its physical form as presented. If, for example, for the purposes of supply or transport, the same chemical is to be presented in a physical form different from that which was tested and which is considered likely to materially alter its performance in a classification test, the substance must also be tested in the new form.

2.12.2.2. A substance or mixture shall be classified as a substance or mixture which in contact with water emits flammable gases if spontaneous ignition takes place in any step of the test procedure.

2.12.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.12.2.

Table 2.12.2

Label elements for substances or mixtures which in contact with water emit flammable gases

Classification	Category 1	Category 2	Category 3
GHS Pictograms			
Signal Word	Danger	Danger	Warning
Hazard Statement	H260: In contact with water releases flammable gases which may ignite spontaneously	H261: In contact with water releases flammable gases	H261: In contact with water releases flammable gases
Precautionary Statement Prevention	P223 P231 + P232 P280	P223 P231 + P232 P280	P231 + P232 P280
Precautionary Statement Response	P335 + P334 P370 + P378	P335 + P334 P370 + P378	P370 + P378
Precautionary Statement Storage	P402 + P404	P402 + P404	P402 + P404

Classification	Category 1	Category 2	Category 3
Precautionary Statement Disposal	P501	P501	P501

2.12.4. ***Additional Classification Considerations***

2.12.4.1. The classification procedure for this class need not be applied if:

(a) the chemical structure of the substance or mixture does not contain metals or metalloids; or

(b) experience in production or handling shows that the substance or mixture does not react with water, e.g. the substance is manufactured with water or washed with water; or

(c) the substance or mixture is known to be soluble in water to form a stable mixture.

2.13. **Oxidising liquids**

2.13.1. ***Definition***

Oxidising liquid means a liquid substance or mixture which, while in itself not necessarily combustible, may, generally by yielding oxygen, cause, or contribute to, the combustion of other material.

2.13.2. ***Classification criteria***

2.13.2.1. An oxidising liquid shall be classified in one of the three categories for this class using test O.2 in Part III, sub-section 34.4.2 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria in accordance with Table 2.13.1:

Table 2.13.1

Criteria for oxidising liquids

Category	Criteria
1	Any substance or mixture which, in the 1:1 mixture, by mass, of substance (or mixture) and cellulose tested, spontaneously ignites; or the mean pressure rise time of a 1:1 mixture, by mass, of substance (or mixture) and cellulose is less than that of a 1:1 mixture, by mass, of 50 % perchloric acid and cellulose.
2	Any substance or mixture which, in the 1:1 mixture, by mass, of substance (or mixture) and cellulose tested, exhibits a mean pressure rise time less than or equal to the mean pressure rise time of a 1:1 mixture, by mass, of 40 % aqueous sodium chlorate solution and cellulose; and the criteria for Category 1 are not met.
3	Any substance or mixture which, in the 1:1 mixture, by mass, of substance (or mixture) and cellulose tested, exhibits a mean pressure rise time less than or equal to the mean pressure rise time of a 1:1 mixture, by mass, of 65 % aqueous nitric acid and cellulose; and the criteria for Category 1 and 2 are not met.

2.13.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.13.2.

Table 2.13.2

Label elements for oxidising liquids

Classification	Category 1	Category 2	Category 3
GHS Pictograms			

Classification	Category 1	Category 2	Category 3
Signal Word	Danger	Danger	Warning
Hazard Statement	H271:May cause fire or explosion; strong oxidiser	H272: May intensify fire; oxidiser	H272: May intensify fire; oxidiser
Precautionary Statement Prevention	P210 P220 P221 P280 P283	P210 P220 P221 P280	P210 P220 P221 P280
Precautionary Statement Response	P306 + P360 P371 + P380 + P375 P370 + P378	P370 + P378	P370 + P378
Precautionary Statement Storage			
Precautionary Statement Disposal	P501	P501	P501

2.13.4. ***Additional Classification Considerations***

2.13.4.1. For organic substances or mixtures the classification procedure for this class shall not apply if:

(a) the substance or mixture does not contain oxygen, fluorine or chlorine; or

(b) the substance or mixture contains oxygen, fluorine or chlorine and these elements are chemically bonded only to carbon or hydrogen.

2.13.4.2. For inorganic substances or mixtures the classification procedure for this class shall not apply if they do not contain oxygen or halogen atoms.

2.13.4.3. In the event of divergence between test results and known experience in the handling and use of substances or mixtures which shows them to be oxidising, judgments based on known experience shall take precedence over test results.

2.13.4.4. In cases where substances or mixtures generate a pressure rise (too high or too low), caused by chemical reactions not characterising the oxidising properties of the substance or mixture, the test described in Part III, sub-section 34.4.2 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria shall be repeated with an inert substance, e.g. diatomite (kieselguhr), in place of the cellulose in order to clarify the nature of the reaction and to check for a false positive result.

2.14. **Oxidising solids**

2.14.1. ***Definition***

Oxidising solid means a solid substance or mixture which, while in itself is not necessarily combustible, may, generally by yielding oxygen, cause, or contribute to, the combustion of other material.

2.14.2. ***Classification criteria***

2.14.2.1. An oxidising solid shall be classified in one of the three categories for this class using test O.1 in Part III, sub-section 34.4.1 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria in accordance with Table 2.14.1:

Table 2.14.1

Criteria for oxidising solids

Category	Criteria
1	Any substance or mixture which, in the 4:1 or 1:1 sample-to-cellulose ratio (by mass) tested, exhibits a mean burning time less than the mean burning time of a 3:2 mixture, by mass, of potassium bromate and cellulose.
2	Any substance or mixture which, in the 4:1 or 1:1 sample-to-cellulose ratio (by mass) tested, exhibits a mean burning time equal to or less than the mean burning time of a 2:3 mixture (by mass) of potassium bromate and cellulose and the criteria for Category 1 are not met.
3	Any substance or mixture which, in the 4:1 or 1:1 sample-to-cellulose ratio (by mass) tested, exhibits a mean burning time equal to or less than the mean burning time of a 3:7 mixture (by mass) of potassium bromate and cellulose and the criteria for Categories 1 and 2 are not met.

Note 1

Some oxidising solids also present explosion hazards under certain conditions (when stored in large quantities). Some types of ammonium nitrate may give rise to an explosion hazard under extreme conditions and the 'Resistance to detonation test' (BC Code, Annex 3, Test 5) can be used to assess this hazard. Appropriate information shall be made in the SDS.

Note 2

The test shall be performed on the substance or mixture in its physical form as presented. If, for example, for the purposes of supply or transport, the same chemical is to be presented in a physical form different from that which was tested and which is considered likely to materially alter its performance in a classification test, the substance shall also be tested in the new form.

2.14.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.14.2.

Table 2.14.2

Label elements for oxidising solids

	Category 1	Category 2	Category 3
GHS Pictograms			
Signal Word	Danger	Danger	Warning
Hazard Statement	H271: May cause fire or explosion; strong oxidiser	H272: May intensify fire; oxidiser	H272: May intensify fire; oxidiser
Precautionary Statement Prevention	P210 P220 P221 P280 P283	P210 P220 P221 P280	P210 P220 P221 P280
Precautionary Statement Response	P306 + P360 P371 + P380 + P375 P370 + P378	P370 + P378	P370 + P378
Precautionary Statement Storage			

	Category 1	Category 2	Category 3
Precautionary Statement Disposal	P501	P501	P501

2.14.4. ***Additional Classification Considerations***

2.14.4.1. For organic substances or mixtures the classification procedure for this class shall not apply if:

(a) the substance or mixture does not contain oxygen, fluorine or chlorine; or

(b) the substance or mixture contains oxygen, fluorine or chlorine and these elements are chemically bonded only to carbon or hydrogen.

2.14.4.2. For inorganic substances or mixtures the classification procedure for this class shall not apply if they do not contain oxygen or halogen atoms.

2.14.4.3. In the event of divergence between test results and known experience in the handling and use of substances or mixtures which shows them to be oxidising, judgments based on known experience shall take precedence over test results.

2.15. **Organic peroxides**

2.15.1. ***Definition***

2.15.1.1. Organic peroxides means liquid or solid organic substances which contain the bivalent -O-O- structure and may be considered derivatives of hydrogen peroxide, where one or both of the hydrogen atoms have been replaced by organic radicals. The term organic peroxide includes organic peroxide mixtures (formulations) containing at least one organic peroxide. Organic peroxides are thermally unstable substances or mixtures, which can undergo exothermic self-accelerating decomposition. In addition, they can have one or more of the following properties:

(i) be liable to explosive decomposition;

(ii) burn rapidly;

(iii) be sensitive to impact or friction;

(iv) react dangerously with other substances.

2.15.1.2. An organic peroxide is regarded as possessing explosive properties when in laboratory testing the mixture (formulation) is liable to detonate, to deflagrate rapidly or to show a violent effect when heated under confinement.

2.15.2. ***Classification criteria***

2.15.2.1. Any organic peroxide shall be considered for classification in this class, unless it contains:

(a) not more than 1,0 % available oxygen from the organic peroxides when containing not more than 1,0 % hydrogen peroxide; or

(b) not more than 0,5 % available oxygen from the organic peroxides when containing more than 1,0 % but not more than 7,0 % hydrogen peroxide.

Note

The available oxygen content (%) of an organic peroxide mixture is given by the formula:

$$16 \times \sum_{i}^{n} \left(\frac{n_i \times c_i}{m_i} \right)$$

where:

n_i = number of peroxygen groups per molecule of organic peroxide i;
c_i = concentration (mass %) of organic peroxide i;
m_i = molecular mass of organic peroxide i.

2.15.2.2. Organic peroxides shall be classified in one of the seven categories of 'Types A to G' for this class, according to the following principles:

(a) any organic peroxide which, as packaged, can detonate or deflagrate rapidly shall be defined as organic peroxide TYPE A;

(b) any organic peroxide possessing explosive properties and which, as packaged, neither detonates nor deflagrates rapidly, but is liable to undergo a thermal explosion in that package shall be defined as organic peroxide TYPE B;

(c) any organic peroxide possessing explosive properties when the substance or mixture as packaged cannot detonate or deflagrate rapidly or undergo a thermal explosion shall be defined as organic peroxide TYPE C;

(d) any organic peroxide which in laboratory testing:

(i) detonates partially, does not deflagrate rapidly and shows no violent effect when heated under confinement; or

(ii) does not detonate at all, deflagrates slowly and shows no violent effect when heated under confinement; or

(iii) does not detonate or deflagrate at all and shows a medium effect when heated under confinement;

shall be defined as organic peroxide TYPE D;

(e) any organic peroxide which, in laboratory testing, neither detonates nor deflagrates at all and shows low or no effect when heated under confinement shall be defined as organic peroxide TYPE E;

(f) any organic peroxide which, in laboratory testing, neither detonates in the cavitated state nor deflagrates at all and shows only a low or no effect when heated under confinement as well as low or no explosive power shall be defined as organic peroxide TYPE F;

(g) any organic peroxide which, in laboratory testing, neither detonates in the cavitated state nor deflagrates at all and shows no effect when heated under confinement nor any explosive power, provided that it is thermally stable, i.e. the SADT is 60 °C or higher for a 50 kg package ([1]), and, for liquid mixtures, a diluent having a boiling point of not less than 150 °C is used for desensitisation, shall be defined as organic peroxide TYPE G. If the organic peroxide is not thermally stable or a diluent having a boiling point less than 150 °C is used for desensitisation, the organic peroxide shall be defined as organic peroxide TYPE F.

Where the test is conducted in the package form and the packaging is changed, a further test shall be conducted where it is considered that the change in packaging will affect the outcome of the test.

2.15.2.3. *Criteria for temperature control*

The following organic peroxides need to be subjected to temperature control:

(a) Organic peroxide types B and C with an SADT ≤ 50 C;

(b) Organic peroxide type D showing a medium effect when heated under confinement ([2]) with an SADT ≤ 50 °C or showing a low or no effect when heated under confinement with an SADT ≤ 45 °C; and

(c) Organic peroxide types E and F with an SADT ≤ 45 °C.

([1]) See UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, sub-sections 28.1, 28.2, 28.3 and Table 28.3.

([2]) As determined by test series E as prescribed in UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, Part II.

Test methods for determining the SADT as well as the derivation of control and emergency temperatures are given in the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, Part II, section 28. The test selected shall be conducted in a manner which is representative, both in size and material, of the package.

2.15.3. ***Hazard Communication***

Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.15.1.

Table 2.15.1

Label elements for organic peroxides

Classification	Type A	Type B	Type C & D	Type E & F	Type G
GHS Pictograms					There are no label elements allocated to this hazard category
Signal Word	Danger	Danger	Danger	Warning	
Hazard Statement	H240: Heating may cause an explosion	H241: Heating may cause a fire or explosion	H242: Heating may cause a fire	H242: Heating may cause a fire	
Precautionary Statement Prevention	P210 P220 P234 P280	P210 P220 P234 P280	P210 P220 P234 P280	P210 P220 P234 P280	
Precautionary Statement Response					
Precautionary Statement Storage	P411 + P235 P410 P420	P411 + P235 P410 P420	P411 + P235 P410 P420	P411 + P235 P410 P420	
Precautionary Statement Disposal	P501	P501	P501	P501	

Type G has no hazard communication elements assigned but shall be considered for properties belonging to other hazard classes.

2.15.4. ***Additional Classification Considerations***

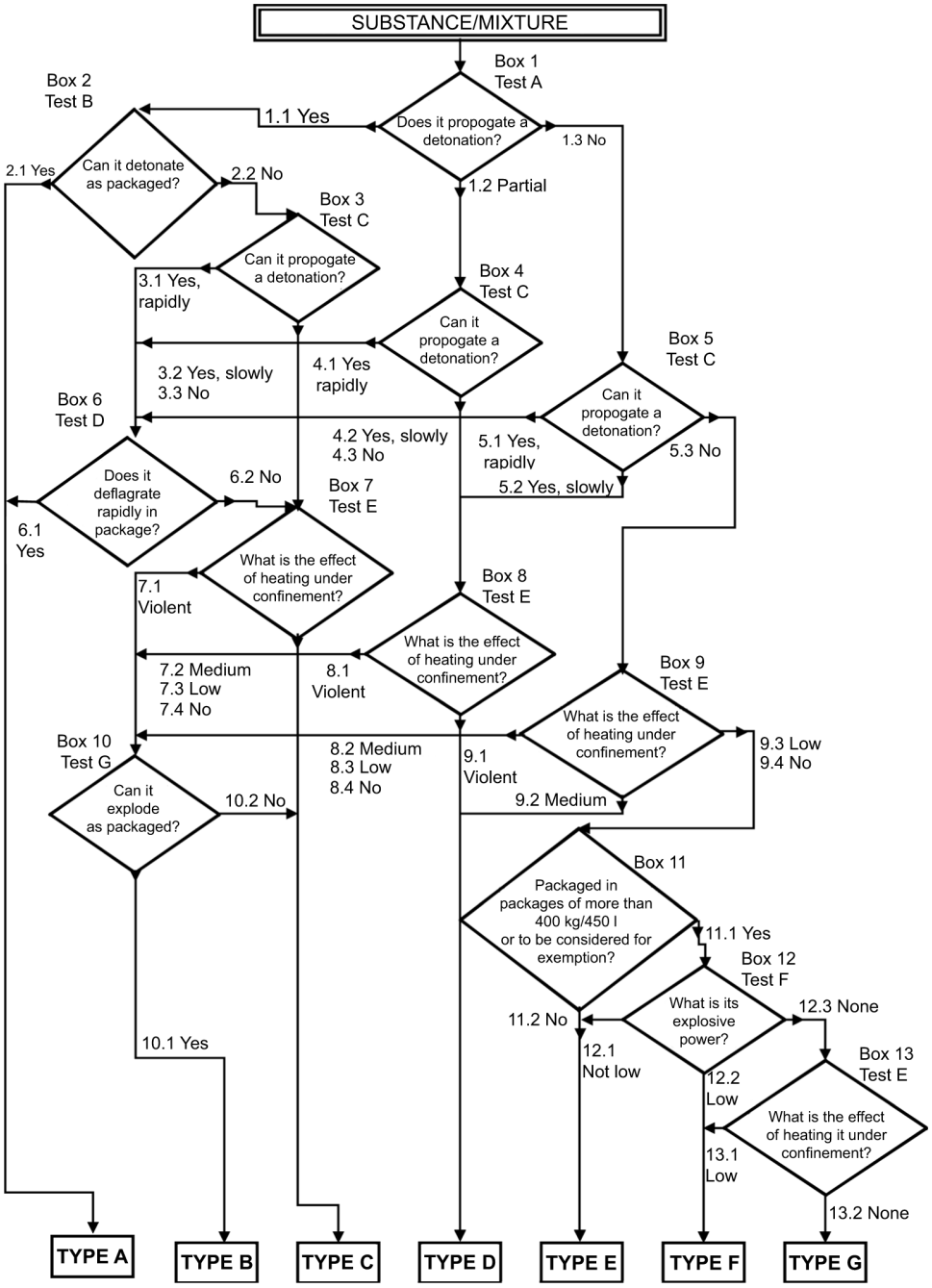
2.15.4.1. Organic peroxides are classified by definition based on their chemical structure and on the available oxygen and hydrogen peroxide contents of the mixture (see 2.15.2.1). The properties of organic peroxides which are necessary for their classification shall be determined experimentally. The classification of organic peroxides shall be performed in accordance with test series A to H as described in Part II of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria. The procedure for classification is described in Figure 2.15.1.

2.15.4.2. Mixtures of already classified organic peroxides may be classified as the same type of organic peroxide as that of the most dangerous component. However, as two stable components can form a thermally less stable mixture, the SADT of the mixture shall be determined.

Note: The sum of the individual parts can be more hazardous than the individual components.

Figure 2.15.1

Organic Peroxides



2.16. **Corrosive to metals**

2.16.1. ***Definition***

A substance or a mixture that is corrosive to metals means a substance or a mixture which by chemical action will materially damage, or even destroy, metals.

2.16.2. ***Classification criteria***

2.16.2.1. A substance or a mixture which is corrosive to metals is classified in a single category for this class, using the test in Part III, sub-section 37.4 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria, in accordance with Table 2.16.1:

Table 2.16.1

Criteria for substances and mixtures corrosive to metals

Category	Criteria
1	Corrosion rate on either steel or aluminium surfaces exceeding 6,25 mm per year at a test temperature of 55 ºC when tested on both materials.

Note

Where an initial test on either steel or aluminium indicates the substance or mixture being tested is corrosive the follow up test on the other metal is not required.

2.16.3. ***Hazard Communication***

Label elements shall be used for substances and mixtures meeting the criteria for classification in this hazard class in accordance with Table 2.16.2.

Table 2.16.2

Label elements for substances and mixtures corrosive to metals

Classification	Category 1
GHS Pictogram	
Signal Word	Warning
Hazard Statement	H290: May be corrosive to metals
Precautionary Statement Prevention	P234
Precautionary Statement Response	P390
Precautionary Statement Storage	P406
Precautionary Statement Disposal	

2.16.4. ***Additional Classification Considerations***

2.16.4.1. The corrosion rate can be measured according to the test method of Part III sub-section 37.4 of the UN Recommendations on the Transport of Dangerous Goods, Manual of Tests and Criteria. The specimen to be used for the test shall be made of the following materials:

(a) for the purposes of testing steel, steel types

— S235JR+CR (1.0037 resp.St 37-2),

— S275J2G3+CR (1.0144 resp.St 44-3), ISO 3574 as amended, Unified Numbering System (UNS) G 10200, or SAE 1020;

(b) for the purposes of testing aluminium: non-clad types 7075-T6 or AZ5GU-T6.

3. PART 3: HEALTH HAZARDS

3.1. **Acute toxicity**

3.1.1. ***Definitions***

3.1.1.1. Acute toxicity means those adverse effects occurring following oral or dermal administration of a single dose of a substance or a mixture, or multiple doses given within 24 hours, or an inhalation exposure of 4 hours.

3.1.1.2. The hazard class Acute Toxicity is differentiated into:

— Acute oral toxicity;

— Acute dermal toxicity;

— Acute inhalation toxicity.

3.1.2. ***Criteria for classification of substances as acutely toxic***

3.1.2.1. Substances can be allocated to one of four toxicity categories based on acute toxicity by the oral, dermal or inhalation route according to the numeric criteria shown in Table 3.1.1. Acute toxicity values are expressed as (approximate) LD_{50} (oral, dermal) or LC_{50} (inhalation) values or as acute toxicity estimates (ATE). Explanatory notes are shown following Table 3.1.1.

Table 3.1.1

Acute toxicity hazard categories and acute toxicity estimates (ATE) defining the respective categories

Exposure Route	Category 1	Category 2	Category 3	Category 4
Oral (mg/kg body-weight) See Note (a)	ATE ≤ 5	5 < ATE ≤ 50	50 < ATE ≤ 300	300 < ATE ≤ 2 000
Dermal (mg/kg bodyweight) See Note (a)	ATE ≤ 50	50 < ATE ≤ 200	200 < ATE ≤ 1 000	1 000 < ATE ≤ 2 000
Gases (ppmV [1]) see: Note (a) Note (b)	ATE ≤ 100	100 < ATE ≤ 500	500 < ATE ≤ 2 500	2 500 < ATE ≤ 20 000
Vapours (mg/l) see: Note (a) Note (b) Note (c)	ATE ≤ 0,5	0,5 < ATE ≤ 2,0	2,0 < ATE ≤ 10,0	10,0 < ATE ≤ 20,0
Dusts and Mists (mg/l) see: Note (a) Note (b)	ATE ≤ 0,05	0,05 < ATE ≤ 0,5	0,5 < ATE ≤ 1,0	1,0 < ATE ≤ 5,0

[1] Gas concentrations are expressed in parts per million per volume (ppmV).

Notes to Table 3.1.1

(a) The acute toxicity estimate (ATE) for the classification of a substance or ingredient in a mixture is derived using:

— the LD_{50}/LC_{50} where available,

— the appropriate conversion value from Table 3.1.2 that relates to the results of a range test, or

— the appropriate conversion value from Table 3.1.2 that relates to a classification category.

(b) Generic concentration limits for inhalation toxicity in the table are based on 4 hour testing exposures. Conversion of existing inhalation toxicity data which have been generated using a 1 hour exposure can be carried out by dividing by a factor of 2 for gases and vapours and 4 for dusts and mists.

(c) For some substances or mixtures the test atmosphere will not just be a vapour but will consist of a mixture of liquid and vapour phases. For other substances or mixtures the test atmosphere may consist of a vapour which is near the gaseous phase. In these latter cases, classification shall be based on ppmV as follows: Category 1 (100 ppmV), Category 2 (500 ppmV), Category 3 (2 500 ppmV), Category 4 (20 000 ppmV).

The terms 'dust', 'mist' and 'vapour' are defined as follows:

— Dust: solid particles of a substance or mixture suspended in a gas (usually air);

— Mist: liquid droplets of a substance or mixture suspended in a gas (usually air);

— Vapour: the gaseous form of a substance or mixture released from its liquid or solid state.

Dust is generally formed by mechanical processes. Mist is generally formed by condensation of supersaturated vapours or by physical shearing of liquids. Dusts and mists generally have sizes ranging from less than 1 to about 100 μm.

3.1.2.2. *Specific considerations for classification of substances as acutely toxic*

3.1.2.2.1. The preferred test species for evaluation of acute toxicity by the oral and inhalation routes is the rat, while the rat or rabbit are preferred for evaluation of acute dermal toxicity. When experimental data for acute toxicity are available in several animal species, scientific judgement shall be used in selecting the most appropriate LD_{50} value from among valid, well-performed tests.

3.1.2.3. *Specific considerations for classification of substances as acutely toxic by the inhalation route*

3.1.2.3.1. Units for inhalation toxicity are a function of the form of the inhaled material. Values for dusts and mists are expressed in mg/l. Values for gases are expressed in ppmV. Acknowledging the difficulties in testing vapours, some of which consist of mixtures of liquid and vapour phases, the table provides values in units of mg/l. However, for those vapours which are near the gaseous phase, classification shall be based on ppmV.

3.1.2.3.2. Of particular importance in classifying for inhalation toxicity is the use of well articulated values in the high toxicity categories for dusts and mists. Inhaled particles between 1 and 4 microns mean mass aerodynamic diameter (MMAD) will deposit in all regions of the rat respiratory tract. This particle size range corresponds to a maximum dose of about 2 mg/l. In order to achieve applicability of animal experiments to human exposure, dusts and mists would ideally be tested in this range in rats.

3.1.2.3.3. In addition to classification for inhalation toxicity, if data are available that indicates that the mechanism of toxicity was corrosivity, the substance or mixture shall also be labelled as 'corrosive to the respiratory tract' (see note 1 in 3.1.4.1). Corrosion of the respiratory tract is defined by destruction of the respiratory tract tissue after a single, limited period of exposure analogous to skin corrosion; this includes destruction of the mucosa. The corrosivity evaluation can be based on expert judgment using such evidence as: human and animal experience, existing (in vitro) data, pH values, information from similar substances or any other pertinent data.

3.1.3. ***Criteria for classification of mixtures as acutely toxic***

3.1.3.1. The criteria for classification of substances for acute toxicity as outlined in section 3.1.2 are based on lethal dose data (tested or derived). For mixtures, it is necessary to obtain or derive information that allows the criteria to be applied to the mixture for the purpose of classification. The approach to classification for acute toxicity is tiered, and is dependent upon the amount of information available for the mixture itself and for its ingredients. The flow chart of Figure 3.1.1 outlines the process to be followed.

3.1.3.2. For acute toxicity each route of exposure shall be considered for the classification of mixtures, but only one route of exposure is needed as long as this route is followed (estimated or tested) for all ingredients. If the acute toxicity is determined for more than one route of exposure, the more severe hazard category will be used for classification. All available information shall be considered and all relevant routes of exposure shall be identified for hazard communication.

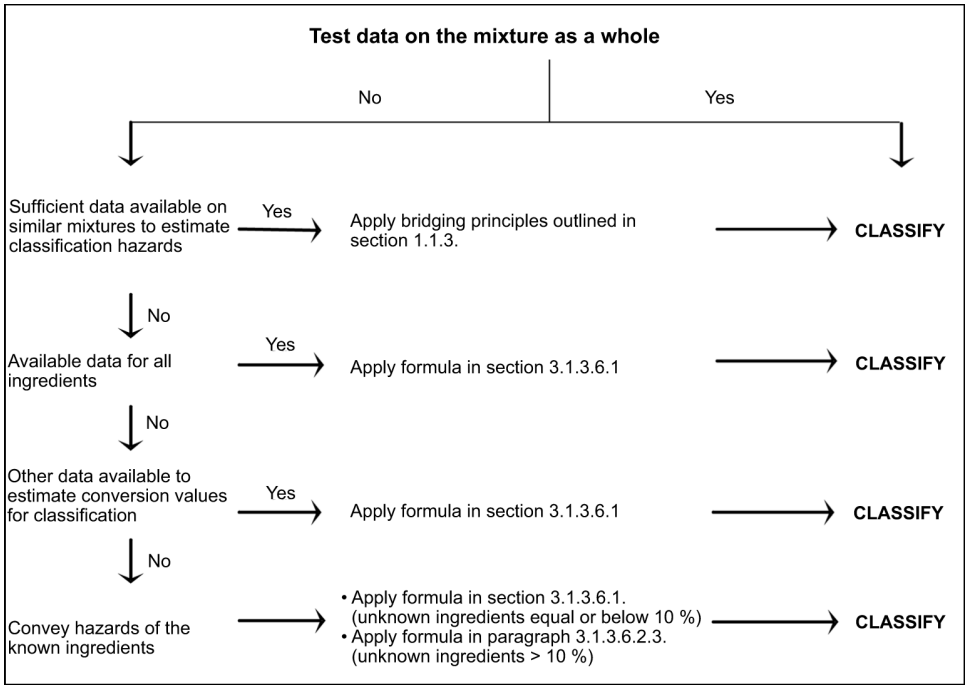
3.1.3.3. In order to make use of all available data for purposes of classifying the hazards of the mixtures, certain assumptions have been made and are applied where appropriate in the tiered approach:

(a) the 'relevant ingredients' of a mixture are those which are present in concentrations of 1 % (w/w for solids, liquids, dusts, mists and vapours and v/v for gases) or greater, unless there is a reason to suspect that an ingredient present at a concentration of less than 1 % is still relevant for classifying the mixture for acute toxicity (see Table 1.1).

(b) where a classified mixture is used as an ingredient of another mixture, the actual or derived acute toxicity estimate (ATE) for that mixture may be used, when calculating the classification of the new mixture using the formulas in section 3.1.3.6.1 and paragraph 3.1.3.6.2.3.

Figure 3.1.1

Tiered approach to classification of mixtures for acute toxicity



3.1.3.4. *Classification of mixtures where acute toxicity data are available for the complete mixture*

3.1.3.4.1. Where the mixture itself has been tested to determine its acute toxicity, it shall be classified according to the same criteria as those used for substances, presented in Table 3.1.1. If test data for the mixture are not available, the procedures presented under sections 3.1.3.5 and 3.1.3.6 shall be followed.

3.1.3.5. *Classification of mixtures where acute toxicity data are available for the complete mixture: bridging principles*

3.1.3.5.1. Where the mixture itself has not been tested to determine its acute toxicity, but there are sufficient data on the individual ingredients and similar tested mixtures to adequately characterise the hazards of the mixture, these data shall be used in accordance with the bridging rules set out in section 1.1.3.

3.1.3.5.2. If a mixture is diluted with water or other totally non-toxic material, the toxicity of the mixture can be calculated from test data on the undiluted mixture.

3.1.3.6. *Classification of mixtures based on ingredients of the mixture (Additivity formula)*

3.1.3.6.1. Data available for all ingredients

In order to ensure that classification of the mixture is accurate, and that the calculation need only be performed once for all systems, sectors, and categories, the acute toxicity estimate (ATE) of ingredients shall be considered as follows:

(a) include ingredients with a known acute toxicity, which fall into any of the acute toxicity categories shown in Table 3.1.1;

(b) ignore ingredients that are presumed not acutely toxic (e.g., water, sugar);

(c) ignore ingredients if the oral limit test does not show acute toxicity at 2 000 mg/kg bodyweight.

Ingredients that fall within the scope of this paragraph are considered to be ingredients with a known acute toxicity estimate (ATE).

The ATE of the mixture is determined by calculation from the ATE values for all relevant ingredients according to the following formula for Oral, Dermal or Inhalation Toxicity:

$$\frac{100}{ATE_{mix}} = \sum_{n} \frac{C_i}{ATE_i}$$

where:

C_i = concentration of ingredient i (% w/w or % v/v)

i = the individual ingredient from 1 to n

n = the number of ingredients

ATE_i = Acute Toxicity Estimate of ingredient i.

3.1.3.6.2. Classification of mixtures when data are not available for all components

3.1.3.6.2.1. Where an ATE is not available for an individual ingredient of the mixture, but available information, such as that listed below, can provide a derived conversion value such as those laid out in Table 3.1.2, the formula in section 3.1.3.6.1 shall be applied.

This includes evaluation of:

(a) extrapolation between oral, dermal and inhalation acute toxicity estimates ([1]). Such an evaluation could require appropriate pharmacodynamic and pharmacokinetic data;

(b) evidence from human exposure that indicates toxic effects but does not provide lethal dose data;

(c) evidence from any other toxicity tests/assays available on the substance that indicates toxic acute effects but does not necessarily provide lethal dose data; or

(d) data from closely analogous substances using structure/activity relationships.

([1]) For ingredients with acute toxicity estimates available for other than the most appropriate exposure route, values may be extrapolated from the available exposure route(s) to the most appropriate route. Dermal and inhalation route data are not always required for ingredients. However, in case data requirements for specific ingredients include acute toxicity estimates for the dermal and inhalation route, the values to be used in the formula need to be from the required exposure route.

This approach generally requires substantial supplemental technical information, and a highly trained and experienced expert (expert judgement, see section 1.1.1), to reliably estimate acute toxicity. If such information is not available, proceed to paragraph 3.1.3.6.2.3.

3.1.3.6.2.2. In the event that an ingredient without any useable information at all is used in a mixture at a concentration of 1 % or greater, it is concluded that the mixture cannot be attributed a definitive acute toxicity estimate. In this situation the mixture shall be classified based on the known ingredients only, with the additional statement that x percent of the mixture consists of ingredient(s) of unknown toxicity.

3.1.3.6.2.3. If the total concentration of the ingredient(s) with unknown acute toxicity is ≤ 10 % then the formula presented in section 3.1.3.6.1 shall be used. If the total concentration of the ingredient(s) with unknown toxicity is > 10 %, the formula presented in section 3.1.3.6.1 shall be corrected to adjust for the total percentage of the unknown ingredient(s) as follows:

$$\frac{100 - \left(\sum C_{\text{unknown if}} > 10\ \%\right)}{ATE_{mix}} = \sum_{n} \frac{C_i}{ATE_i}$$

Table 3.1.2

Conversion from experimentally obtained acute toxicity range values (or acute toxicity hazard categories) to acute toxicity point estimates for classification for the respective routes of exposure

Exposure routes	Classification Category or experimentally obtained acute toxicity range estimate	Converted acute toxicity point estimate (see Note 1)
Oral (mg/kg bodyweight)	0 < Category 1 ≤ 5	0,5
	5 < Category 2 ≤ 50	5
	50 < Category 3 ≤ 300	100
	300 < Category 4 ≤ 2 000	500
Dermal (mg/kg bodyweight)	0 < Category 1 ≤ 50	5
	50 < Category 2 ≤ 200	50
	200 < Category 3 ≤ 1 000	300
	1 000 < Category 4 ≤ 2 000	1 100
Gases (ppmV)	0 < Category 1 ≤ 100	10
	100 < Category 2 ≤ 500	100
	500 < Category 3 ≤ 2 500	700
	2 500 < Category 4 ≤ 20 000	4 500
Vapours (mg/l)	0 < Category 1 ≤ 0,5	0,05
	0,5 < Category 2 ≤ 2,0	0,5
	2,0 < Category 3 ≤ 10,0	3
	10,0 < Category 4 ≤ 20,0	11
Dust/mist (mg/l)	0< Category 1 ≤ 0,05	0,005
	0,05 < Category 2 ≤ 0,5	0,05
	0,5 < Category 3 ≤ 1,0	0,5
	1,0 < Category 4 ≤ 5,0	1,5

Note 1

These values are designed to be used in the calculation of the ATE for classification of a mixture based on its components and do not represent test results.

3.1.4. ***Hazard Communication***

3.1.4.1. Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 3.1.3.

Table 3.1.3

Acute toxicity label elements

Classification	Category 1	Category 2	Category 3	Category 4
GHS Pictograms				
Signal Word	Danger	Danger	Danger	Warning
Hazard Statement: — Oral	H300: Fatal if swallowed	H300: Fatal if swallowed	H301: Toxic if swallowed	H302: Harmful if swallowed
— Dermal	H310:Fatal in contact with skin	H310:Fatal in contact with skin	H311: Toxic in contact with skin	H312: Harmful in contact with skin
— Inhalation (see Note 1)	H330:Fatal if inhaled	H330: Fatal if inhaled	H331: Toxic if inhaled	H332: Harmful if inhaled
Precautionary Statement Prevention (oral)	P264 P270	P264 P270	P264 P270	P264 P270
Precautionary Statement Response (oral)	P301 + P310 P321 P330	P301 + P310 P321 P330	P301 + P310 P321 P330	P301 + P312 P330
Precautionary Statement Storage (oral)	P405	P405	P405	
Precautionary Statement Disposal (oral)	P501	P501	P501	P501
Precautionary Statement Prevention (dermal)	P262 P264 P270 P280	P262 P264 P270 P280	P280	P280
Precautionary Statement Response (dermal)	P302 + P350 P310 P322 P361 P363	P302 + P350 P310 P322 P361 P363	P302 + P352 P312 P322 P361 P363	P302 + P352 P312 P322 P363
Precautionary Statement Storage (dermal)	P405	P405	P405	
Precautionary Statement Disposal (dermal)	P501	P501	P501	P501
Precautionary Statement Prevention (inhalation)	P260 P271 P284	P260 P271 P284	P261 P271	P261 P271
Precautionary Statement Response (inhalation)	P304 + P340 P310 P320	P304 + P340 P310 P320	P304 + P340 P311 P321	P304 + P340 P312
Precautionary Statement Storage (inhalation)	P403 + P233 P405	P403 + P233 P405	P403 + P233 P405	
Precautionary Statement Disposal (inhalation)	P501	P501	P501	

Note 1

In addition to classification for inhalation toxicity, if data are available that indicates that the mechanism of toxicity is corrosivity, the substance or mixture shall also be labelled as EUH071: 'corrosive to the respiratory tract' — see advice at 3.1.2.3.3. In addition to an appropriate acute toxicity pictogram, a corrosivity pictogram (used for skin and eye corrosivity) may be added together with the statement 'corrosive to the respiratory tract'.

Note 2

In the event that an ingredient without any useable information at all is used in a mixture at a concentration of 1 % or greater, the mixture shall be labelled with the additional statement that 'x percent of the mixture consists of ingredient(s) of unknown toxicity' — see advice at 3.1.3.6.2.2.

3.2. **Skin corrosion/irritation**

3.2.1. ***Definitions***

3.2.1.1. Skin Corrosion means the production of irreversible damage to the skin; namely, visible necrosis through the epidermis and into the dermis, following the application of a test substance for up to 4 hours. Corrosive reactions are typified by ulcers, bleeding, bloody scabs, and, by the end of observation at 14 days, by discolouration due to blanching of the skin, complete areas of alopecia, and scars. Histopathology shall be considered to evaluate questionable lesions.

Skin Irritation means the production of reversible damage to the skin following the application of a test substance for up to 4 hours.

3.2.2. ***Classification criteria for substances***

3.2.2.1. Several factors need to be considered in determining the corrosion and irritation potential of substances before testing is undertaken. Solid substances (powders) may become corrosive or irritant when moistened or in contact with moist skin or mucous membranes. Existing human experience and animal data from single or repeated exposure shall be the first line of analysis, as they give information directly relevant to effects on the skin. In vitro alternatives that have been validated and accepted may also be used to help make classification decisions (see Article 5). In some cases enough information may be available from structurally related compounds to make classification decisions.

3.2.2.2. Likewise, pH extremes like ≤ 2 and ≥ 11,5 may indicate the potential to cause skin effects, especially when buffering capacity is known, although the correlation is not perfect. Generally, such substances are expected to produce significant effects on the skin. If consideration of alkali/acid reserve suggests the substance may not be corrosive despite the low or high pH value, then further testing shall be carried out to confirm this, preferably by use of an appropriate validated in vitro test.

3.2.2.3. If a substance is highly toxic by the dermal route, a skin irritation/corrosion study is not practicable since the amount of test substance to be applied considerably exceeds the toxic dose and, consequently, results in the death of the animals. When observations are made of skin irritation/corrosion in acute toxicity studies and are observed up through the limit dose, additional testing is not needed, provided that the dilutions used and species tested are equivalent.

3.2.2.4. All the above information that is available on a substance shall be used in determining the need for in vivo skin irritation testing.

Although information might be gained from the evaluation of single parameters within a tier (see paragraph 3.2.2.5), e.g. caustic alkalis with extreme pH shall be considered as skin corrosives, there is merit in considering the totality of existing information and making an overall weight of evidence determination. This is especially true when there is information available on some but not all parameters. Generally, primary emphasis shall be placed upon existing human experience and data, followed by animal experience and testing data, followed by other sources of information, but case-by-case determinations are necessary.

3.2.2.5. A tiered approach to the evaluation of initial information shall be considered, where applicable, recognising that all elements may not be relevant in certain cases.

3.2.2.6. *Corrosion*

3.2.2.6.1. On the basis of the results of animal testing a substance is classified as corrosive, as shown in Table 3.2.1. A corrosive substance is a substance that produces destruction of skin tissue, namely, visible necrosis through the epidermis and into the dermis, in at least 1 tested animal after exposure up to a 4 hour duration. Corrosive reactions are typified by ulcers, bleeding, bloody scabs and, by the end of observation at 14 days, by discoloration due to blanching of the skin, complete areas of alopecia and scars. Histopathology shall be considered to discern questionable lesions.

3.2.2.6.2. Three subcategories are provided within the corrosive category: subcategory 1A — where responses are noted following up to 3 minutes exposure and up to 1 hour observation; subcategory 1B — where responses are described following exposure between 3 minutes and 1 hour and observations up to 14 days; and subcategory 1C — where responses occur after exposures between 1 hour and 4 hours and observations up to 14 days.

3.2.2.6.3. The use of human data is discussed in paragraphs 3.2.2.1 and 3.2.2.4 and also in paragraphs 1.1.1.3, 1.1.1.4 and.1.1.1.5.

Table 3.2.1

Skin Corrosive category and subcategories

		Corrosive in > 1 of 3 animals	
	Corrosive subcategories	Exposure	Observation
Category 1: Corrosive	1A	≤ 3 minutes	≤ 1 hour
	1B	> 3 minutes - ≤ 1 hour	≤ 14 days
	1C	> 1 hour - ≤ 4 hours	≤ 14 days

3.2.2.7. *Irritation*

3.2.2.7.1. Using the results of animal testing a single irritant category (Category 2) is presented in Table 3.2.2. The use of human data is discussed in paragraphs 3.2.2.1 and 3.2.2.4 and also in paragraphs 1.1.1.3, 1.1.1.4 and.1.1.1.5. The major criterion for the irritant category is that at least 2 of 3 tested animals have a mean score of ≥ 2,3 - ≤ 4,0.

Table 3.2.2

Skin irritation category

Category	Criteria
Category 2: Irritant	(1) Mean value of ≥ 2,3 - ≤ 4,0 for erythema/eschar or for oedema in at least 2 of 3 tested animals from gradings at 24, 48 and 72 hours after patch removal or, if reactions are delayed, from grades on 3 consecutive days after the onset of skin reactions; or (2) Inflammation that persists to the end of the observation period normally 14 days in at least 2 animals, particularly taking into account alopecia (limited area), hyperkeratosis, hyperplasia, and scaling; or (3) In some cases where there is pronounced variability of response among animals, with very definite positive effects related to chemical exposure in a single animal but less than the criteria above.

3.2.2.8. *Comments on responses obtained in skin irritation tests in animals*

3.2.2.8.1. Animal irritant responses within a test can be quite variable, as they are with corrosion. The major criterion for classification of a substance as irritant to skin, as shown in paragraph 3.2.2.7.1, is the mean value of the scores for either erythema/eschar or oedema calculated in at least 2 of 3 tested animals. A separate irritant criterion accommodates cases when there is a significant irritant response but less than the mean score criterion for a positive test. For example, a test material might be designated as an irritant if at least 1 of 3 tested animals shows a very elevated mean score throughout the study, including lesions persisting at the end of an observation period of normally 14 days. Other responses could also fulfil this criterion. However, it should be ascertained that the responses are the result of chemical exposure.

3.2.2.8.2. Reversibility of skin lesions is another consideration in evaluating irritant responses. When inflammation persists to the end of the observation period in 2 or more test animals, taking into consideration alopecia (limited area), hyperkeratosis, hyperplasia and scaling, then a material shall be considered to be an irritant.

3.2.3. ***Classification criteria for mixtures***

3.2.3.1. *Classification of mixtures when data are available for the complete mixture*

3.2.3.1.1. The mixture will be classified using the criteria for substances, and taking into account the testing and evaluation strategies to develop data for these hazard classes.

3.2.3.1.2. Unlike other hazard classes, there are alternative tests available for skin corrosivity of certain types of substances and mixtures that can give an accurate result for classification purposes, as well as being simple and relatively inexpensive to perform. When considering testing of the mixture, classifiers are encouraged to use a tiered weight of evidence strategy as included in the criteria for classification of substances for skin corrosion and irritation (paragraph 3.2.2.5), to help ensure an accurate classification as well as avoid unnecessary animal testing. A mixture is considered corrosive to skin (Skin Corrosive Category 1) if it has a pH of 2 or less or a pH of 11,5 or greater. If consideration of alkali/acid reserve suggests the substance or mixture may not be corrosive despite the low or high pH value, then further testing shall be carried out to confirm this, preferably by use of an appropriate validated in vitro test.

3.2.3.2. *Classification of mixtures when data are not available for the complete mixture: bridging principles*

3.2.3.2.1. Where the mixture itself has not been tested to determine its skin irritation/corrosion hazards, but there are sufficient data on the individual ingredients and similar tested mixtures to adequately characterise the hazards of the mixture, these data shall be used in accordance with the bridging rules set out in section 1.1.3.

3.2.3.3. *Classification of mixtures when data are available for all components or only for some components of the mixture*

3.2.3.3.1. In order to make use of all available data for purposes of classifying the skin irritation/corrosion hazards of mixtures, the following assumption has been made and is applied where appropriate in the tiered approach:

Assumption: the 'relevant ingredients' of a mixture are those which are present in concentrations of 1 % (w/w for solids, liquids, dusts, mists and vapours and v/v for gases) or greater, unless there is a presumption (e.g., in the case of corrosive ingredients) that an ingredient present at a concentration of less than 1 % can still be relevant for classifying the mixture for skin irritation/corrosion.

3.2.3.3.2. In general, the approach to classification of mixtures as irritant or corrosive to skin when data are available on the components, but not on the mixture as a whole, is based on the theory of additivity, such that each corrosive or irritant component contributes to the overall irritant or corrosive properties of the mixture in proportion to its potency and concentration. A weighting factor of 10 is used for corrosive components when they are present at a concentration below the generic concentration limit for classification with Category 1, but are at a concentration that will contribute to the classification of the mixture as an irritant. The mixture is classified as corrosive or irritant when the sum of the concentrations of such components exceeds a concentration limit.

3.2.3.3.3. Table 3.2.3 provides the generic concentration limits to be used to determine if the mixture is considered to be an irritant or a corrosive to the skin.

3.2.3.3.4.1. Particular care must be taken when classifying certain types of mixtures containing substances such as acids and bases, inorganic salts, aldehydes, phenols, and surfactants. The approach explained in paragraphs 3.2.3.3.1 and 3.2.3.3.2 may not be applicable given that many of such substances are corrosive or irritant at concentrations < 1 %.

3.2.3.3.4.2. For mixtures containing strong acids or bases the pH shall be used as a classification criterion (see paragraph 3.2.3.1.2) since pH is a better indicator of corrosion than the concentration limits of Table 3.2.3.

3.2.3.3.4.3. A mixture containing ingredients that are corrosive or irritant to the skin and that cannot be classified on the basis of the additivity approach (Table 3.2.3), due to chemical characteristics that make this approach unworkable, shall be classified as Skin Corrosive Category 1A, 1B or 1C if it contains ≥ 1 % of an ingredient classified in Category 1A, 1B or 1C respectively or as Category 2 when it contains ≥ 3 % of an irritant ingredient. Classification of mixtures with ingredients for which the approach in Table 3.2.3 does not apply is summarised in Table 3.2.4.

3.2.3.3.5. On occasion, reliable data may show that the skin corrosion/irritation hazard of an ingredient will not be evident when present at a level above the generic concentration limits mentioned in Tables 3.2.3 and 3.2.4. In these cases the mixture shall be classified according to that data (see also Articles 10 and 11). On other occasions, when it is expected that the skin corrosion/irritation hazard of an ingredient is not evident when present at a level above the generic concentration limits mentioned in Tables 3.2.3 and 3.2.4, testing of the mixture shall be considered. In those cases the tiered weight of evidence strategy shall be applied, as described in paragraph 3.2.2.5.

3.2.3.3.6. If there are data showing that (an) ingredient(s) is/are corrosive or irritant at a concentration of < 1 % (corrosive) or < 3 % (irritant), the mixture shall be classified accordingly.

Table 3.2.3

Generic concentration limits of ingredients classified for skin corrosive/irritant hazard (Category 1 or 2) that trigger classification of the mixture as corrosive/irritant to skin

Sum of ingredients classified as:	Concentration triggering classification of a mixture as:	
	Skin Corrosive	Skin Irritant
	Category 1 (see note below)	Category 2
Skin Corrosive Categories 1A, 1B, 1C	≥ 5 %	≥ 1 % but < 5 %
Skin irritant Category 2		≥ 10 %
(10 × Skin Corrosive Category 1A, 1B, 1C) + Skin irritant Category 2		≥ 10 %

Note

The sum of all ingredients of a mixture classified as Skin Corrosive Category 1A, 1B or 1C respectively, shall each be ≥ 5 % respectively in order to classify the mixture as either Skin Corrosive Category 1A, 1B or 1C. If the sum of the Skin Corrosive Category 1A ingredients is < 5 % but the sum of Category 1A+1B ingredients is ≥ 5 %, the mixture shall be classified as Skin Corrosive Category 1B. Similarly, if the sum of Skin Corrosive Category 1A+1B ingredients is < 5 % but the sum of Category 1A+1B+1C ingredients is ≥ 5 % the mixture shall be classified as Skin Corrosive Category 1C.

Table 3.2.4

Generic concentration limits of ingredients of a mixture for which the additivity approach does not apply, that trigger classification of the mixture as corrosive/irritant to skin

Ingredient:	Concentration:	Mixture classified as: Skin
Acid with pH ≤ 2	≥ 1 %	Category 1
Base with pH ≥ 11,5	≥ 1 %	Category 1
Other corrosive (Categories 1A, 1B, 1C) ingredients for which additivity does not apply	≥ 1 %	Category 1
Other irritant (Category 2) ingredients for which additivity does not apply, including acids and bases	≥ 3 %	Category 2

3.2.4. ***Hazard Communication***

3.2.4.1. Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 3.2.5.

Table 3.2.5

Label elements for skin corrosion/irritation

Classification	Category 1 A/1 B/1 C	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H314: Causes severe skin burns and eye damage	H315: Causes skin irritation
Precautionary Statement Prevention	P260 P264 P280	P264 P280
Precautionary Statement Response	P301 + P330 + P331 P303 + P361 + P353 P363 P304 + P340 P310 P321 P305 + P351 + P338	P302 + P352 P321 P332 + P313 P362
Precautionary Statement Storage	P405	
Precautionary Statement Disposal	P501	

3.3. **Serious eye damage/eye irritation**

3.3.1. ***Definitions***

3.3.1.1. Serious eye damage means the production of tissue damage in the eye, or serious physical decay of vision, following application of a test substance to the anterior surface of the eye, which is not fully reversible within 21 days of application.

Eye irritation means the production of changes in the eye following the application of test substance to the anterior surface of the eye, which are fully reversible within 21 days of application.

3.3.2. ***Classification criteria for substances***

3.3.2.1. The classification system for substances involves a tiered testing and evaluation scheme, combining pre-existing information on serious ocular tissue damage and on eye irritation (including data relating to historical human or animal experience) as well as considerations on (Q)SAR and the output of validated in vitro tests in order to avoid unnecessary animal testing.

3.3.2.2. Before any in vivo testing for serious eye damage/eye irritation is carried out, all existing information on a substance shall be reviewed. Preliminary decisions can often be made from existing data as to whether a substance causes serious (i.e. irreversible) damage to the eyes. If a substance can be classified on the basis of these data, no testing is required.

3.3.2.3. Several factors need to be considered in determining the serious eye damage or irritation potential of a substance before testing is undertaken. Accumulated human and animal experience shall be the first line of analysis, as it gives information directly relevant to effects on the eye. In some cases enough information may be available from structurally related compounds to make hazard decisions. Likewise, pH extremes like ≤ 2 and ≥ 11,5 may produce serious eye damage, especially when associated with significant buffering capacity. Such substances are expected to produce significant effects on the eyes. Possible skin corrosion has to be

evaluated prior to consideration of serious eye damage/eye irritation in order to avoid testing for local effects on eyes with skin corrosive substances. Skin corrosive substances shall be considered as leading to serious damage to the eyes as well (Category 1), while skin irritant substances may be considered as leading to eye irritation (Category 2). In vitro alternatives that have been validated and accepted can be used to make classification decisions (see Article 5).

3.3.2.4. All the above information that is available on a substance shall be used in determining the need for in vivo eye irritation testing. Although information may be gained from the evaluation of single parameters within a tier (e.g. caustic alkalis with extreme pH shall be considered as local corrosives), the totality of existing information shall be considered in making an overall weight of evidence determination, particularly when there is information available on some but not all parameters. Generally, primary emphasis shall be placed upon expert judgement, considering human experience with the substance, followed by the outcome of skin irritation testing and of well-validated alternative methods. Animal testing with corrosive substances or mixtures shall be avoided whenever possible.

3.3.2.5. A tiered approach to the evaluation of initial information shall be considered where applicable, while recognising that all elements may not be relevant in certain cases.

3.3.2.6. *Irreversible effects on the eye/serious damage to eyes (Category 1)*

3.3.2.6.1. Substances that have the potential to seriously damage the eyes are classified in Category 1 (irreversible effects on the eye). Substances are classified in this hazard category on the basis of the results of animal testing, in accordance with the criteria listed in Table 3.3.1. These observations include animals with grade 4 cornea lesions and other severe reactions (e.g., destruction of cornea) observed at any time during the test, as well as persistent corneal opacity, discoloration of the cornea by a dye substance, adhesion, pannus, and interference with the function of the iris or other effects that impair sight. In this context, persistent lesions are considered those which are not fully reversible within an observation period of normally 21 days. Substances are also classified in Category 1 if they fulfil the criteria of corneal opacity ≥ 3 or iritis > 1,5 detected in a Draize eye test with rabbits, recognising that such severe lesions usually do not reverse within a 21-day observation period.

Table 3.3.1

Category for irreversible eye effects

Category	Criteria
Irreversible effects on the eye (Category 1)	If, when applied to the eye of an animal, a substance produces: — at least in one animal effects on the cornea, iris or conjunctiva that are not expected to reverse or have not fully reversed within an observation period of normally 21 days; and/or — at least in 2 of 3 tested animals, a positive response of: — corneal opacity ≥ 3 and/or — iritis > 1,5 calculated as the mean scores following grading at 24, 48 and 72 hours after installation of the test material.

3.3.2.6.2. The use of human data is discussed in paragraphs 3.3.2.1, 3.3.2.4, and also in paragraphs 1.1.1.3, 1.1.1.4 and 1.1.1.5.

3.3.2.7. *Reversible effects on the eye (Category 2)*

3.3.2.7.1. Substances that have the potential to induce reversible eye irritation are classified in Category 2 (irritating to eyes).

Table 3.3.2

Category for reversible eye effects

Category	Criteria
Irritating to eyes (Category 2)	if, when applied to the eye of an animal, a substance produces: — at least in 2 of 3 tested animals, a positive response of: — corneal opacity ≥ 1 and/or — iritis ≥ 1, and/or — conjunctival redness ≥ 2 and/or — conjunctival oedema (chemosis) ≥ 2 — calculated as the mean scores following grading at 24, 48 and 72 hours after installation of the test material, and which fully reverses within an observation period of 21 days

3.3.2.7.2. For those substances where there is pronounced variability among animal responses, this information shall be taken into account in determining the classification.

3.3.3. ***Classification criteria for mixtures***

3.3.3.1. *Classification of mixtures when data are available for the complete mixture*

3.3.3.1.1. The mixture will be classified using the criteria for substances, and taking into account the testing and evaluation strategies used to develop data for these hazard classes.

3.3.3.1.2. Unlike other hazard classes, there are alternative tests available for skin corrosivity of certain types of mixtures that give an accurate result for classification purposes, as well as being simple and relatively inexpensive to perform. When considering testing of the mixture classifiers are encouraged to use a tiered weight of evidence strategy as included in the criteria for classification of substances for skin corrosion and serious eye damage and eye irritation to help ensure an accurate classification, as well as avoid unnecessary animal testing. A mixture is considered to cause serious eye damage (Category 1) if it has a pH ≤ 2,0 or ≥ 11,5. If consideration of alkali/acid reserve suggests the mixture may not have the potential to cause serious eye damage despite the low or high pH value, then further testing needs to be carried out to confirm this, preferably by use of an appropriate validated in vitro test.

3.3.3.2. *Classification of mixtures when data are not available for the complete mixture: bridging principles*

3.3.3.2.1. Where the mixture itself has not been tested to determine its skin corrosivity or potential to cause serious eye damage or irritation, but there are sufficient data on the individual ingredients and similar tested mixtures to adequately characterise the hazards of the mixture, these data shall be used in accordance with the bridging rules set out in section 1.1.3.

3.3.3.3. *Classification of mixtures when data are available for all components or only for some components of the mixture*

3.3.3.3.1. In order to make use of all available data for purposes of classifying the eye irritation/serious eye damaging properties of the mixtures, the following assumption has been made and is applied where appropriate in the tiered approach:

Assumption: The 'relevant ingredients' of a mixture are those which are present in concentrations of 1 % (w/w for solids, liquids, dusts, mists and vapours and v/v for gases) or greater, unless there is a presumption (e.g. in the case of corrosive ingredients) that an ingredient present at a concentration of less than 1 % is still relevant for classifying the mixture for eye irritation/serious eye damage.

3.3.3.3.2. In general, the approach to classification of mixtures as eye irritant or seriously damaging to the eye when data are available on the components, but not on the mixture as a whole, is based on the theory of additivity, such that each corrosive or irritant component contributes to the overall irritant or corrosive properties of the mixture in proportion to its potency and concentration. A weighting factor of 10 is used for corrosive components when they are present at a concentration below the generic concentration limit for classification in Category 1, but are at a concentration that will contribute to the classification of the mixture as an irritant. The mixture is classified as seriously damaging to the eye or eye irritant when the sum of the concentrations of such components exceeds a concentration limit.

3.3.3.3.3. Table 3.3.3 provides the generic concentration limits to be used to determine if the mixture shall be classified as irritant or as seriously damaging to the eye.

3.3.3.3.4.1. Particular care must be taken when classifying certain types of mixtures containing substances such as acids and bases, inorganic salts, aldehydes, phenols, and surfactants. The approach explained in paragraphs 3.3.3.3.1 and 3.3.3.3.2 might not work given that many of such substances are corrosive or irritant at concentrations < 1 %.

3.3.3.3.4.2. For mixtures containing strong acids or bases the pH shall be used as classification criteria (see paragraph 3.3.2.3) since pH will be a better indicator of serious eye damage than the generic concentration limits of Table 3.3.3.

3.3.3.3.4.3. A mixture containing corrosive or irritant ingredients that cannot be classified based on the additivity approach (Table 3.3.3), due to chemical characteristics that make this approach unworkable, shall be classified as Category 1 for effects on the eye if it contains ≥ 1 % of a corrosive ingredient and as Category 2 when it contains ≥ 3 % of an irritant ingredient. Classification of mixtures with ingredients for which the approach in Table 3.3.3 does not apply is summarised in Table 3.3.4.

3.3.3.3.5. On occasion, reliable data may show that the reversible/irreversible eye effects of an ingredient will not be evident when present at a level above the generic concentration limits mentioned in Tables 3.3.3 and 3.3.4. In these cases the mixture shall be classified according to those data. On other occasions, when it is expected that the skin corrosion/irritation hazards or the reversible/irreversible eye effects of an ingredient will not be evident when present at a level above the generic concentration limits mentioned in Tables 3.3.3 and 3.3.4, testing of the mixture shall be considered. In those cases, the tiered weight of evidence strategy shall be applied.

3.3.3.3.6. If there are data showing that (an) ingredient(s) may be corrosive or irritant at a concentration of < 1 % (corrosive) or < 3 % (irritant), the mixture shall be classified accordingly.

Table 3.3.3

Generic concentration limits of ingredients of a mixture classified as Skin corrosive Category 1 and/or eye Category 1 or 2 for effects on the eye that trigger classification of the mixture for effects on the eye (Category 1 or 2)

Sum of ingredients classified as:	Concentration triggering classification of a mixture as:	
	Irreversible Eye Effects	Reversible Eye Effects
	Category 1	Category 2
Eye Effects Category 1 or Skin Corrosive Category 1A, 1B, 1C	≥ 3 %	≥ 1 % but < 3 %
Eye Effects Category 2		≥ 10 %
(10 × Eye Effects Category 1) + Eye effects Category 2		≥ 10 %
Skin Corrosive Category 1A, 1B, 1C + Eye effects Category 1	≥ 3 %	≥ 1 % but < 3 %
10 × (Skin Corrosive Category 1A, 1B, 1C + Eye Effects Category 1) + Eye Effects Category 2		≥ 10 %

Table 3.3.4

Generic concentration limits of ingredients of a mixture for which the additivity approach does not apply, that trigger classification of the mixture as hazardous to the eye

Ingredient	Concentration	Mixture classified as: Eye
Acid with pH ≤ 2	≥ 1 %	Category 1
Base with pH ≥ 11,5	≥ 1 %	Category 1
Other corrosive (Category 1) ingredients for which additivity does not apply	≥ 1 %	Category 1

Ingredient	Concentration	Mixture classified as: Eye
Other irritant (Category 2) ingredients for which additivity does not apply, including acids and bases	≥ 3 %	Category 2

3.3.4. ***Hazard Communication***

3.3.4.1. Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 3.3.5.

Table 3.3.5

Label elements for serious eye damage/eye irritation

Classification	Category 1	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H318: Causes serious eye damage	H319: Causes serious eye irritation
Precautionary Statement Prevention	P280	P264 P280
Precautionary Statement Response	P305 + P351 + P338 P310	P305 + P351 + P338 P337 + P313
Precautionary Statement Storage		
Precautionary Statement Disposal		

3.4. **Respiratory or skin sensitisation**

3.4.1. ***Definitions and general considerations***

3.4.1.1. Respiratory sensitiser means a substance that will lead to hypersensitivity of the airways following inhalation of the substance.

3.4.1.2. Skin sensitiser means a substance that will lead to an allergic response following skin contact.

3.4.1.3. For the purpose of section 3.4, sensitisation includes two phases: the first phase is induction of specialised immunological memory in an individual by exposure to an allergen. The second phase is elicitation, i.e. production of a cell-mediated or antibody-mediated allergic response by exposure of a sensitised individual to an allergen.

3.4.1.4. For respiratory sensitisation, the pattern of induction followed by elicitation phases is shared in common with skin sensitisation. For skin sensitisation, an induction phase is required in which the immune system learns to react; clinical symptoms can then arise when subsequent exposure is sufficient to elicit a visible skin reaction (elicitation phase). As a consequence, predictive tests usually follow this pattern in which there is an induction phase, the response to which is measured by a standardised elicitation phase, typically involving a patch test. The local lymph node assay is the exception, directly measuring the induction response. Evidence of skin sensitisation in humans normally is assessed by a diagnostic patch test.

3.4.1.5. Usually, for both skin and respiratory sensitisation, lower levels are necessary for elicitation than are required for induction. Provisions for alerting sensitised individuals to the presence of a particular sensitiser in a mixture can be found at section 3.4.4.

3.4.1.6. The hazard class Respiratory or Skin Sensitisation is differentiated into:

— Respiratory Sensitisation;

— Skin Sensitisation.

3.4.2. ***Classification criteria for substances***

3.4.2.1. *Respiratory sensitisers*

Substances shall be classified as respiratory sensitisers (Category 1) in accordance with the criteria in Table 3.4.1:

Table 3.4.1

Hazard category for respiratory sensitisers

Category	Criteria
Category 1	Substances shall be classified as respiratory sensitisers (Category 1) in accordance with the following criteria: (a) if there is evidence in humans that the substance can lead to specific respiratory hypersensitivity and/or (b) if there are positive results from an appropriate animal test.

3.4.2.1.1 Human evidence

3.4.2.1.1.1. Evidence that a substance can induce specific respiratory hypersensitivity will normally be based on human experience. In this context, hypersensitivity is normally seen as asthma, but other hypersensitivity reactions such as rhinitis/conjunctivitis and alveolitis are also considered. The condition will have the clinical character of an allergic reaction. However, immunological mechanisms do not have to be demonstrated.

3.4.2.1.1.2. When considering the human evidence, it is necessary for a decision on classification to take into account, in addition to the evidence from the cases:

(a) the size of the population exposed;

(b) the extent of exposure.

The use of human data is discussed in paragraphs 1.1.1.3, 1.1.1.4 and 1.1.1.5.

3.4.2.1.1.3. The evidence referred to above could be

(a) clinical history and data from appropriate lung function tests related to exposure to the substance, confirmed by other supportive evidence which may include:

(i) in vivo immunological test (e.g. skin prick test);

(ii) in vitro immunological test (e.g. serological analysis);

(iii) studies that indicate other specific hypersensitivity reactions where immunological mechanisms of action have not been proven, e.g. repeated low-level irritation, pharmacologically mediated effects;

(iv) a chemical structure related to substances known to cause respiratory hypersensitivity;

(b) data from one or more positive bronchial challenge tests with the substance conducted according to accepted guidelines for the determination of a specific hypersensitivity reaction.

3.4.2.1.1.4. Clinical history shall include both medical and occupational history to determine a relationship between exposure to a specific substance and development of respiratory hypersensitivity. Relevant information includes aggravating factors both in the home and workplace, the onset and progress of the disease, family history and medical history of the patient in question. The medical history shall also include a note of other allergic or airway disorders from childhood, and smoking history.

3.4.2.1.1.5. The results of positive bronchial challenge tests are considered to provide sufficient evidence for classification on their own. It is however recognised that in practice many of the examinations listed above will already have been carried out.

3.4.2.1.2. Animal studies

3.4.2.1.2.1. Data from appropriate animal studies [1] which may be indicative of the potential of a substance to cause sensitisation by inhalation in humans [2] may include:

(i) measurements of Immunoglobulin E (IgE) and other specific immunological parameters in mice;

(ii) specific pulmonary responses in guinea pigs.

3.4.2.2. *Skin sensitisers*

3.4.2.2.1. Substances shall be classified as skin sensitisers (Category 1) in accordance with the criteria in Table 3.4.2:

Table 3.4.2

Hazard category for skin sensitisers

Category	Criteria
Category 1	Substances shall be classified as skin sensitisers (Category 1) in accordance with the following criteria: (i) if there is evidence in humans that the substance can lead to sensitisation by skin contact in a substantial number of persons, or (ii) if there are positive results from an appropriate animal test (see specific criteria in paragraph 3.4.2.2.4.1).

3.4.2.2.2. Specific considerations

3.4.2.2.2.1. For classification of a substance as a skin sensitiser, evidence shall include any or all of the following:

(a) positive data from patch testing, normally obtained in more than one dermatology clinic;

(b) epidemiological studies showing allergic contact dermatitis caused by the substance; Situations in which a high proportion of those exposed exhibit characteristic symptoms are to be looked at with special concern, even if the number of cases is small;

(c) positive data from appropriate animal studies;

(d) positive data from experimental studies on humans (see Article 7(3));

(e) well documented episodes of allergic contact dermatitis, normally obtained in more than one dermatology clinic.

The use of human data is discussed in paragraphs 1.1.1.3, 1.1.1.4 and 1.1.1.5.

3.4.2.2.2.2. Positive effects seen in either humans or animals will normally justify classification. Evidence from animal studies (see section 3.4.2.2.4) is usually much more reliable than evidence from human exposure. However, in cases where evidence is available from both sources, and there is conflict between the results, the quality and reliability of the evidence from both sources must be assessed in order to resolve the question of classification on a case-by-case basis. Normally, human data are not generated in controlled experiments with volunteers for the purpose of hazard classification but rather as part of risk assessment to confirm lack of effects seen in animal tests. Consequently, positive human data on skin sensitisation are usually derived from case-control or other, less defined studies. Evaluation of human data must therefore be carried out with caution, as the frequency of cases reflect, in addition to the intrinsic properties of the substances, factors such as the exposure situation, bioavailability, individual predisposition and preventive measures taken. Negative human data can not normally be used to negate positive results from animal studies.

[1] At present recognised animal models for the testing of respiratory hypersensitivity are not available.

[2] The mechanisms by which substances induce symptoms of asthma are not yet fully known. For preventative measures, these substances are considered respiratory sensitisers. However, if on the basis of the evidence, it can be demonstrated that these substances induce symptoms of asthma by irritation only in people with bronchial hyper reactivity, they should not be considered as respiratory sensitisers.

3.4.2.2.2.3. If none of the above mentioned conditions are met the substance need not be classified as a skin sensitiser. However, a combination of two or more indicators of skin sensitisation as listed below may alter the decision, which shall be considered on a case-by-case basis:

(a) isolated episodes of allergic contact dermatitis;

(b) epidemiological studies of limited power, e.g. where chance, bias or confounders have not been ruled out fully with reasonable confidence;

(c) data from animal tests, performed according to existing guidelines, which do not meet the criteria for a positive result described in paragraph 3.4.2.2.4.1, but which are sufficiently close to the limit to be considered significant;

(d) positive data from non-standard methods;

(e) positive results from close structural analogues.

3.4.2.2.3. Immunological contact urticaria

3.4.2.2.3.1. Some substances meeting the criteria for classification as respiratory sensitisers may in addition cause immunological contact urticaria. Consideration shall be given to classifying these substances also as skin sensitisers and including information concerning contact urticaria on the label or in the SDS using appropriate warning information.

3.4.2.2.3.2. For substances which produce signs of immunological contact urticaria but which do not fulfil the criteria as a respiratory sensitiser, consideration shall be given to classification as a skin sensitiser. There is no recognised animal model available to identify substances which cause immunological contact urticaria. Therefore, classification will normally be based on human evidence, which will be similar to that for skin sensitisation.

3.4.2.2.4. Animal studies

3.4.2.2.4.1. When an adjuvant type guinea pig test method for skin sensitisation is used, a response of at least 30 % of the animals is considered as positive. For a non-adjuvant guinea pig test method a response of at least 15 % of the animals is considered positive. Test methods for skin sensitisation described in Regulation (EC) No 440/2008 adopted in accordance with Article 13(3) of Regulation (EC) No 1907/2006 ('Test Method Regulation') shall be used, or other methods provided that they are well-validated and scientific justification is given.

3.4.3. ***Classification criteria for mixtures***

3.4.3.1. *Classification of mixtures when data are available for the complete mixture*

3.4.3.1.1. When reliable and good quality evidence from human experience or appropriate studies in experimental animals, as described in the criteria for substances, is available for the mixture, then the mixture can be classified by weight of evidence evaluation of these data. Care shall be exercised in evaluating data on mixtures, that the dose used does not render the results inconclusive.

3.4.3.2. *Classification of mixtures when data are not available for the complete mixture: bridging principles*

3.4.3.2.1. Where the mixture itself has not been tested to determine its sensitising properties, but there are sufficient data on the individual ingredients and similar tested mixtures to adequately characterise the hazards of the mixture, these data shall be used in accordance with the bridging rules set out in section 1.1.3.

3.4.3.3. *Classification of mixtures when data are available for all ingredients or only for some ingredients of the mixture*

3.4.3.3.1. The mixture shall be classified as a respiratory or skin sensitiser when at least one ingredient has been classified as a respiratory or skin sensitiser and is present at or above the appropriate generic concentration limit as shown in Table 3.4.3 for solid/liquid and gas respectively.

3.4.3.3.2. Some substances that are classified as sensitisers may elicit a response, when present in a mixture in quantities below the concentrations established in Table 3.4.1, in individuals who are already sensitised to the substance or mixture (see Note 1 to Table 3.4.3).

Table 3.4.3

Generic concentration limits of ingredients of a mixture classified as either skin sensitisers or respiratory sensitisers that trigger classification of the mixture

Ingredient classified as:	Concentration triggering classification of a mixture as:		
	Skin Sensitiser	Respiratory Sensitiser	
	All physical states	Solid/Liquid	Gas
Skin Sensitiser	≥ 0,1 % (Note 1)	—	—
	≥ 1,0 % (Note 2)	—	—
Respiratory Sensitiser	—	≥ 0,1 % (Note 1)	≥ 0,1 % (Note 1)
	—	≥ 1,0 % (Note 3)	≥ 0,2 % (Note 3)

Note 1

This concentration limit is generally used for the application of the special labelling requirements of Annex II section 2.8 to protect already sensitised individuals. A SDS is required for the mixture containing an ingredient above this concentration.

Note 2

This concentration limit is used to trigger classification of a mixture as a skin sensitiser.

Note 3

This concentration limit is used to trigger classification of a mixture as a respiratory sensitiser.

3.4.4. ***Hazard communication***

3.4.4.1. Label elements shall be used for substances or mixtures meeting the criteria for classification in this hazard class in accordance with Table 3.4.4.

Table 3.4.4

Respiratory or skin sensitisation label elements.

Classification	Respiratory sensitisation	Skin sensitisation
	Category 1	Category 1
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H334: May cause allergy or asthma symptoms or breathing difficulties if inhaled	H317: May cause an allergic skin reaction
Precautionary Statement Prevention	P261 P285	P261 P272 P280

Classification	Respiratory sensitisation	Skin sensitisation
	Category 1	Category 1
Precautionary Statement Response	P304 + P341 P342+ P311	P302 + P352 P333 + P313 P321 P363
Precautionary Statement Storage		
Precautionary Statement Disposal	P501	P501

3.5. **Germ cell mutagenicity**

3.5.1. ***Definitions and general considerations***

3.5.1.1. A mutation means a permanent change in the amount or structure of the genetic material in a cell. The term 'mutation' applies both to heritable genetic changes that may be manifested at the phenotypic level and to the underlying DNA modifications when known (including specific base pair changes and chromosomal translocations). The term 'mutagenic' and 'mutagen' will be used for agents giving rise to an increased occurrence of mutations in populations of cells and/or organisms.

3.5.1.2. The more general terms 'genotoxic' and 'genotoxicity' apply to agents or processes which alter the structure, information content, or segregation of DNA, including those which cause DNA damage by interfering with normal replication processes, or which in a non-physiological manner (temporarily) alter its replication. Genotoxicity test results are usually taken as indicators for mutagenic effects.

3.5.2. ***Classification criteria for substances***

3.5.2.1. This hazard class is primarily concerned with substances that may cause mutations in the germ cells of humans that can be transmitted to the progeny. However, the results from mutagenicity or genotoxicity tests in vitro and in mammalian somatic and germ cells in vivo are also considered in classifying substances and mixtures within this hazard class.

3.5.2.2. For the purpose of classification for germ cell mutagenicity, substances are allocated to one of two categories as shown in Table 3.5.1.

Table 3.5.1

Hazard categories for germ cell mutagens

Categories	Criteria
CATEGORY 1:	Substances known to induce heritable mutations or to be regarded as if they induce heritable mutations in the germ cells of humans. Substances known to induce heritable mutations in the germ cells of humans.
Category 1A:	The classification in Category 1A is based on positive evidence from human epidemiological studies. Substances to be regarded as if they induce heritable mutations in the germ cells of humans.

Categories	Criteria
Category 1B:	The classification in Category 1B is based on: — positive result(s) from in vivo heritable germ cell mutagenicity tests in mammals; or — positive result(s) from in vivo somatic cell mutagenicity tests in mammals, in combination with some evidence that the substance has potential to cause mutations to germ cells. It is possible to derive this supporting evidence from mutagenicity/genotoxicity tests in germ cells in vivo, or by demonstrating the ability of the substance or its metabolite(s) to interact with the genetic material of germ cells; or — positive results from tests showing mutagenic effects in the germ cells of humans, without demonstration of transmission to progeny; for example, an increase in the frequency of aneuploidy in sperm cells of exposed people.
CATEGORY 2:	Substances which cause concern for humans owing to the possibility that they may induce heritable mutations in the germ cells of humans The classification in Category 2 is based on: — positive evidence obtained from experiments in mammals and/or in some cases from in vitro experiments, obtained from: — somatic cell mutagenicity tests in vivo, in mammals; or — other in vivo somatic cell genotoxicity tests which are supported by positive results from in vitro mutagenicity assays. Note: Substances which are positive in in vitro mammalian mutagenicity assays, and which also show chemical structure activity relationship to known germ cell mutagens, shall be considered for classification as Category 2 mutagens.

3.5.2.3. *Specific considerations for classification of substances as germ cell mutagens*

3.5.2.3.1. To arrive at a classification, test results are considered from experiments determining mutagenic and/or genotoxic effects in germ and/or somatic cells of exposed animals. Mutagenic and/or genotoxic effects determined in in vitro tests shall also be considered.

3.5.2.3.2. The system is hazard based, classifying substances on the basis of their intrinsic ability to induce mutations in germ cells. The scheme is, therefore, not meant for the (quantitative) risk assessment of substances.

3.5.2.3.3. Classification for heritable effects in human germ cells is made on the basis of well conducted, sufficiently validated tests, preferably as described in Regulation (EC) No 440/2008 adopted in accordance with Article 13(3) of Regulation (EC) No 1907/2006 ('Test Method Regulation') such as those listed in the following paragraphs. Evaluation of the test results shall be done using expert judgement and all the available evidence shall be weighed in arriving at a classification.

3.5.2.3.4. In vivo heritable germ cell mutagenicity tests, such as:

— rodent dominant lethal mutation test;

— mouse heritable translocation assay.

3.5.2.3.5. In vivo somatic cell mutagenicicty tests, such as:

— mammalian bone marrow chromosome aberration test;

— mouse spot test;

— mammalian erythrocyte micronucleus test.

3.5.2.3.6. Mutagenicity/genotoxicity tests in germ cells, such as:

(a) mutagenicity tests:

— mammalian spermatogonial chromosome aberration test;

— spermatid micronucleus assay;

(b) Genotoxicity tests:

— sister chromatid exchange analysis in spermatogonia;

— unscheduled DNA synthesis test (UDS) in testicular cells.

3.5.2.3.7. Genotoxicity tests in somatic cells such as:

— liver Unscheduled synthesis test (UDS) in vivo;

— mammalian bone marrow Sister Chromatid Exchanges (SCE);

3.5.2.3.8. In vitro mutagenicity tests such as:

— in vitro mammalian chromosome aberration test;

— in vitro mammalian cell gene mutation test;

— bacterial reverse mutation tests.

3.5.2.3.9. The classification of individual substances shall be based on the total weight of evidence available, using expert judgement (See 1.1.1). In those instances where a single well-conducted test is used for classification, it shall provide clear and unambiguously positive results. If new, well validated, tests arise these may also be used in the total weight of evidence to be considered. The relevance of the route of exposure used in the study of the substance compared to the route of human exposure shall also be taken into account.

3.5.3. ***Classification criteria for mixtures***

3.5.3.1. *Classification of mixtures when data are available for all ingredients or only for some ingredients of the mixture*

3.5.3.1.1. The mixture shall be classified as a mutagen when at least one ingredient has been classified as a Category 1A, Category 1B or Category 2 mutagen and is present at or above the appropriate generic concentration limit as shown in Table 3.5.2 for Category 1A, Category 1B and Category 2 respectively.

Table 3.5.2

Generic concentration limits of ingredients of a mixture classified as germ cell mutagens that trigger classification of the mixture.

Ingredient classified as:	Concentration limits triggering classification of a mixture as:		
	Category 1A mutagen	Category 1B mutagen	Category 2 mutagen
Category 1A mutagen	≥ 0,1 %	—	—
Category 1B mutagen	—	≥ 0,1 %	—
Category 2 mutagen	—	—	≥ 1,0 %

Note

The concentration limits in the table above apply to solids and liquids (w/w units) as well as gases (v/v units).

3.5.3.2. *Classification of mixtures when data are available for the complete mixture*

3.5.3.2.1. Classification of mixtures will be based on the available test data for the individual ingredients of the mixture using concentration limits for the ingredients classified as germ cell mutagens. On a case-by-case basis, test data on mixtures may be used for classification when demonstrating effects that have not been established from the evaluation based on the individual ingredients. In such cases, the test results for the mixture as a whole must be shown to be conclusive taking into account dose and other factors such as duration, observations, sensitivity and statistical analysis of germ cell mutagenicity test systems. Adequate documentation supporting the classification shall be retained and made available for review upon request.

3.5.3.3. *Classification of mixtures when data are not available for the complete mixture: bridging principles*

3.5.3.3.1. Where the mixture itself has not been tested to determine its germ cell mutagenicity hazard, but there are sufficient data on the individual ingredients and similar tested mixtures (subject to paragraph 3.5.3.2.1), to adequately characterise the hazards of the mixture, these data shall be used in accordance with the applicable bridging rules set out in section 1.1.3.

3.5.4. ***Hazard communication***

3.5.4.1. Label elements shall be used in accordance with Table 3.5.3, for substances or mixtures meeting the criteria for classification in this hazard class.

Table 3.5.3

Label elements of germ cell mutagenicity

Classification	Category 1A or Category 1B	Category 2
GHS Pictograms		
Signal Word	Danger	Warning
Hazard Statement	H340: May cause genetic defects (state route of exposure if it is conclusively proven that no other routes of exposure cause the hazard)	H341: Suspected of causing genetic defects (state route of exposure if it is conclusively proven that no other routes of exposure cause the hazard)
Precautionary Statement Prevention	P201 P202 P281	P201 P202 P281
Precautionary Statement Response	P308 + P313	P308 + P313
Precautionary Statement Storage	P405	P405
Precautionary Statement Disposal	P501	P501

3.5.5. ***Additional classification considerations***

It is increasingly accepted that the process of chemical-induced tumorigenesis in humans and animals involves genetic changes for example in proto-oncogenes and/or tumour suppresser genes of somatic cells. Therefore, the demonstration of mutagenic properties of substances in somatic and/or germ cells of mammals in vivo may have implications for the potential classification of these substances as carcinogens (see also Carcinogenicity, section 3.6, paragraph 3.6.2.2.6).

3.6. **Carcinogenicity**

3.6.1. ***Definition***

3.6.1.1. Carcinogen means a substance or a mixture of substances which induce cancer or increase its incidence. Substances which have induced benign and malignant tumours in well performed experimental studies on animals are considered also to be presumed or suspected human carcinogens unless there is strong evidence that the mechanism of tumour formation is not relevant for humans.

3.6.2. ***Classification criteria for substances***

3.6.2.1. For the purpose of classification for carcinogenicity, substances are allocated to one of two categories based on strength of evidence and additional considerations (weight of evidence). In certain instances, route-specific classification may be warranted, if it can be conclusively proved that no other route of exposure exhibits the hazard.

Table 3.6.1

Hazard categories for carcinogens

Categories	Criteria
CATEGORY 1:	Known or presumed human carcinogens A substance is classified in Category 1 for carcinogenicity on the basis of epidemiological and/or animal data. A substance may be further distinguished as:
Category 1A:	Category 1A, known to have carcinogenic potential for humans, classification is largely based on human evidence, or
Category 1B:	Category 1B, presumed to have carcinogenic potential for humans, classification is largely based on animal evidence. The classification in Category 1A and 1B is based on strength of evidence together with additional considerations (see section 3.6.2.2). Such evidence may be derived from: — human studies that establish a causal relationship between human exposure to a substance and the development of cancer (known human carcinogen); or — animal experiments for which there is sufficient [1] evidence to demonstrate animal carcinogenicity (presumed human carcinogen). In addition, on a case-by-case basis, scientific judgement may warrant a decision of presumed human carcinogenicity derived from studies showing limited evidence of carcinogenicity in humans together with limited evidence of carcinogenicity in experimental animals.
CATEGORY 2:	Suspected human carcinogens The placing of a substance in Category 2 is done on the basis of evidence obtained from human and/or animal studies, but which is not sufficiently convincing to place the substance in Category 1A or 1B, based on strength of evidence together with additional considerations (see section 3.6.2.2). Such evidence may be derived either from limited [1] evidence of carcinogenicity in human studies or from limited evidence of carcinogenicity in animal studies.

[1] Note: See 3.6.2.2.4.

3.6.2.2. *Specific considerations for classification of substances as carcinogens*

3.6.2.2.1. Classification as a carcinogen is made on the basis of evidence from reliable and acceptable studies and is intended to be used for substances which have an intrinsic property to cause cancer. The evaluations shall be based on all existing data, peer-reviewed published studies and additional acceptable data.

3.6.2.2.2. Classification of a substance as a carcinogen is a process that involves two interrelated determinations: evaluations of strength of evidence and consideration of all other relevant information to place substances with human cancer potential into hazard categories.

3.6.2.2.3. Strength of evidence involves the enumeration of tumours in human and animal studies and determination of their level of statistical significance. Sufficient human evidence demonstrates causality between human exposure and the development of cancer, whereas sufficient evidence in animals shows a causal relationship between the substance and an increased incidence of tumours. Limited evidence in humans is demonstrated by a positive association between exposure and cancer, but a causal relationship cannot be stated. Limited evidence in animals is provided when data suggest a carcinogenic effect, but are less than sufficient. The terms 'sufficient' and 'limited' have been used here as they have been defined by the International Agency for Research on Cancer (IARC) and read as follows:

(a) Carcinogenicity in humans

The evidence relevant to carcinogenicity from studies in humans is classified into one of the following categories:

— sufficient evidence of carcinogenicity: a causal relationship has been established between exposure to the agent and human cancer. That is, a positive relationship has been observed between the exposure and cancer in studies in which chance, bias and confounding could be ruled out with reasonable confidence;

— limited evidence of carcinogenicity: a positive association has been observed between exposure to the agent and cancer for which a causal interpretation is considered to be credible, but chance, bias or confounding could not be ruled out with reasonable confidence.

(b) Carcinogenicity in experimental animals

Carcinogenicity in experimental animals can be evaluated using conventional bioassays, bioassays that employ genetically modified animals, and other in-vivo bioassays that focus on one or more of the critical stages of carcinogenesis. In the absence of data from conventional long-term bioassays or from assays with neoplasia as the end-point, consistently positive results in several models that address several stages in the multistage process of carcinogenesis should be considered in evaluating the degree of evidence of carcinogenicity in experimental animals. The evidence relevant to carcinogenicity in experimental animals is classified into one of the following categories:

— sufficient evidence of carcinogenicity: a causal relationship has been established between the agent and an increased incidence of malignant neoplasms or of an appropriate combination of benign and malignant neoplasms in (a) two or more species of animals or (b) two or more independent studies in one species carried out at different times or in different laboratories or under different protocols. An increased incidence of tumours in both sexes of a single species in a well-conducted study, ideally conducted under Good Laboratory Practices, can also provide sufficient evidence. A single study in one species and sex might be considered to provide sufficient evidence of carcinogenicity when malignant neoplasms occur to an unusual degree with regard to incidence, site, type of tumour or age at onset, or when there are strong findings of tumours at multiple sites;

— limited evidence of carcinogenicity: the data suggest a carcinogenic effect but are limited for making a definitive evaluation because, e.g. (a) the evidence of carcinogenicity is restricted to a single experiment; (b) there are unresolved questions regarding the adequacy of the design, conduct or interpretation of the studies; (c) the agent increases the incidence only of benign neoplasms or lesions of uncertain neoplastic potential; or (d) the evidence of carcinogenicity is restricted to studies that demonstrate only promoting activity in a narrow range of tissues or organs.

3.6.2.2.4. Additional considerations (as part of the weight of evidence approach (see 1.1.1)). Beyond the determination of the strength of evidence for carcinogenicity, a number of other factors need to be considered that influence the overall likelihood that a substance poses a carcinogenic hazard in humans. The full list of factors that influence this determination would be very lengthy, but some of the more important ones are considered here.

3.6.2.2.5. The factors can be viewed as either increasing or decreasing the level of concern for human carcinogenicity. The relative emphasis accorded to each factor depends upon the amount and coherence of evidence bearing on each. Generally aaieLc)

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-448-00-1	2,3,5,6-tetrafluorobenzoic acid	416-800-1	652-18-6	Skin Irrit. 2 Eye Dam. 1	H315 H318	GHS05 Dgr	H315 H318			
607-449-00-7	reaction mass of: 4,4',4"-[(2,4,6-trioxo-1,3,5(2*H*,4*H*,6*H*)-triazine-1,3,5-triyl)tris[methylene(3,5,5-trimethyl-3,1-cyclohexanediyl)iminocarbonyloxy-2,1-ethanediyl(ethyl)amino]]trisbenzenediazoniumtri[bis(2-methylpropyl)naphthalenesulfonate]; 4,4',4",4"'-[[5,5'-[carbonylbis[imino(1,5,5-trimethyl-3,1-cyclohexanediyl)methylene]]-2,4,6-trioxo-1,3,5(2*H*,4*H*,6*H*)-triazine-1,1',3,3'-tetrayl]tetrakis[methylene(3,5,5-trimethyl-3,1-cyclohexanediyl)iminocarbonyloxy-2,1-ethanediyl(ethyl)amino]]tetrakisbenzenediazoniumtetra[bis(2-methylpropyl)naphthalenesulfonate]	417-080-1	—	Self-react. D **** Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H242 H317 H400 H410	GHS02 GHS07 GHS09 Dgr	H242 H317 H410			
607-450-00-2	2-mercaptobenzothiazolyl-(Z)-(2-aminothiazol-4-yl)-2-(*tert*-butoxycarbonyl) isopropoxyiminoacetate	419-040-9	89604-92-2	Aquatic Chronic 4	H413	—	H413			
607-451-00-8	4-[4-amino-5-hydroxy-3-(4-(2-sulfoxyethylsulfonyl)phenylazo)-2,7-disulfonapht-6-ylazo]-6-[3-(4-amino-5-hydroxy-3-(4-(2-sulfoxyethylsulfonyl)phenylazo)-2,7-disulfonapht-6-ylazo]phenylcarbonylamino]benzenesulfonic acid, sodium salt	417-640-5	161935-19-9	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
607-453-00-9	4-benzyl-2,6-dihydroxy-4-aza-heptylene bis (2,2-dimethyloctanoate)	418-100-1	172964-15-7	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
607-454-00-4	reaction mass of: *trans*-2-(1-methylethyl)-1,3-dioxane-5-carboxylic acid; *cis*-2-(1-methylethyl)-1,3-dioxane-5-carboxylic acid	418-170-3	116193-72-7	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			
607-455-00-X	1-amino-4-(3-[4-chloro-6-(2,5-di-sulfophenylamino)-1,3,5-triazin-2-ylamino]-2,2-dimethyl-propylamino)-anthraquinone-2-sulfonic acid, sodium/lithium salt	419-520-8	172890-93-6	Skin Sens. 1	H317	GHS07 Wng	H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-456-00-5	3-amino-4-chlorobenzoic acid, hexadecyl ester	419-700-6	143269-74-3	Aquatic Chronic 2	H411	GHS09	H411			
607-457-00-0	tetrasodium dihydrogen 1,1''-dihydroxy-8,8''-[p-phenylbis(imino-{6-[4-(2-aminoethyl)piperazin-1-yl]}-1,3,5-triazine-4,2-diyl-imino)]bis(2,2'-azonaphthalene-1',3,6-trisulfonate)	420-350-1	172277-97-3	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			
607-458-00-6	reaction mass of: 2-ethyl-[2,6-dibromo-4-[1-[3,5-dibromo-4-(2-hydroxyethoxy)phenyl]-1-methylethyl]phenoxy]propenoate; 2,2'-diethyl-[4,4'-bis(2,6-dibromophenoxy)-1-methylethylidene] dipropenoate; 2,2'-[(1-methylethylidene)bis[[2,6-dibromo-4,1-phenylene)oxy]ethanol]]	420-850-1	—	Aquatic Chronic 2	H411	GHS09	H411			
607-459-00-1	isopentyl 4-{2-[5-cyano-1,2,3,6-tetrahydro-1-(2-isopropoxyethoxy-carbonylmethyl)-4-methyl-2,6-dioxo-3-pyridylidene]hydrazino} benzoate	418-930-4	—	Aquatic Chronic 4	H413	—	H413			
607-460-00-7	3-tridecyloxy-propyl-ammonium 9-octadecenoate	418-990-1	778577-53-0	STOT RE 2 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H319 H315 H400 H410	GHS08 GHS07 GHS09 Wng	H373 ** H319 H315 H410			
607-461-00-2	reaction mass of: pentasodium 2-{4-{3-methyl-4-[6-sulfonato-4-(2-sulfonato-phenylazo)-naphthalen-1-ylazo]-phenylamino}-6-[3-(2-sulfato-ethanesulfonyl)-phenylamino]-1,3,5-triazin-2-ylamino}-benzene-1,4-disulfonate; pentasodium 2-{4-{3-methyl-4-[7-sulfonato-4-(2-sulfonato-phenylazo)-naphthalen-1-ylazo]-phenylamino}-6-[3-(2-sulfato-ethanesulfonyl)-phenylamino]-1,3,5-triazin-2-ylamino}-benzene-1,4-disulfonate	421-160-1	—	Aquatic Chronic 3	H412	—	H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-462-00-8	reaction mass of: 1-hexyl acetate; 2-methyl-1-pentyl acetate; 3-methyl-1-pentyl acetate; 4-methyl-1-pentyl acetate; other mixed linear and branched C_6-alkyl acetates	421-230-1	88230-35-7	Aquatic Chronic 2	H411	GHS09	H411			
607-463-00-3	3-(phenothiazin-10-yl)propionic acid	421-260-5	362-03-8	Aquatic Chronic 2	H411	GHS09	H411			
607-464-00-9	reaction mass of: 7-chloro-1-ethyl-6-fluoro-1,4-dihydro-4-oxo-quinoline-3-carboxylic acid; 5-chloro-1-ethyl-6-fluoro-1,4-dihydro-4-oxo-quinoline-3-carboxylic acid	421-280-4		Aquatic Chronic 3	H412	—	H412			
607-465-00-4	tris(2-hydroxyethyl)ammonium 7-{4-[4-(2-cyanoamino-4-hydroxy-6-oxidopyrimidin-5-ylazo)benzamido]-2-ethoxy-phenylazo} naphthalene-1,3-disulfonate	421-440-3	778583-04-3	Aquatic Chronic 3	H412	—	H412			
607-466-00-X	reaction mass of: phenyl 1-(1-[2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl]-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate; phenyl 2-(1-(2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl)-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate; phenyl 3-(1-(2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl)-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate	421-480-1	—	Aquatic Chronic 2	H411	GHS09	H411			
607-467-00-5	1,1,3,3-tetrabutyl-1,3-ditinoxydicaprylate	419-430-9	56533-00-7	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H373 ** H314 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H312 H302 H373 ** H314 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-468-00-0	reaction mass of: monosodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; disodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; trisodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; tetrasodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate	419-450-8	—	Skin Sens. 1	H317	GHS07 Wng	H317			
607-469-00-6	disodium 7-((4,6-bis(3-diethylaminopropylamino)-1,3,5-triazine-2-yl)amino)-4-hydroxy-3-(4-(4-sulfonatophenylazo)phenylazo)-2-naphthalene sulfonate	419-460-2	120029-06-3	Aquatic Chronic 3	H412	—	H412			
607-470-00-1	potassium sodium 6,13-dichloro-3,10-bis{2-[4-[3-(2-hydroxysulphonyloxyethanesulfonyl)phenylamino]-6-(2,5-disulfonatophenylamino)-1,3,5-triazin-2-ylamino]ethylamino}benzo[5,6][1,4]oxazino[2,3-b]phenoxazine-4,11-disulfonate	414-100-0	154336-20-6	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			
607-472-00-2	ammonium iron(III) trimethylenediaminetetraacetate hemihydrate	400-660-3	111687-36-6	Aquatic Chronic 2	H411	GHS09	H411			
607-474-00-3	(4-(4-(4-dimethylaminobenzyliden-1-yl)-3-methyl-5-oxo-2-pyrazolin-1-yl)benzoic acid	410-430-4	117573-89-4	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-475-00-9	reaction mass of: tetrasodium 7-(4-[4-chloro-6-[methyl-(3-sulfonatophenyl) amino]-1,3,5-triazin-2-ylamino]-2-ureidophenylazo)naphthalene-1,3,6-trisulfonate; tetrasodium 7-(4-[4-chloro-6-[methyl-(4-sulfonatophenyl)amino]-1,3,5-triazin-2-ylamino]-2-ureidophenylazo)naphthalene-1,3,6-trisulfonate (1:1)	412-940-2	148878-18-6	Skin Sens. 1	H317	GHS07 Wng	H317			
607-476-00-4	trisodium *N*,*N*-bis(carboxymethyl)-β-alanine	414-070-9	129050-62-0	Skin Corr. 1B Aquatic Chronic 3	H314 H412	GHS05 Dgr	H314 H412			
607-478-00-5	tetramethylammonium hydrogen phthalate	416-900-5	79723-02-7	Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1	H301 H373 ** H400	GHS06 GHS08 GHS09 Dgr	H301 H373 ** H400			
607-479-00-0	hexadecyl 4-chloro-3-[2-(5,5-dimethyl-2,4-dioxo-1,3-oxazolidin-3-yl)-4,4-dimethyl-3-oxopentamido]benzoate	418-550-9	168689-49-4	Aquatic Chronic 4	H413	—	H413			
607-480-00-6	1,2-benzenedicarboxylic acid; di-C_{7-11}-branched and linear alkylesters	271-084-6	68515-42-4	Repr. 1B	H360-Df	GHS08 Dgr	H360Df			
607-487-00-4	reaction mass of: disodium 4-(3-ethoxycarbonyl-4-(5-(3-ethoxycarbonyl-5-hydroxy-1-(4-sulfonatophenyl)pyrazol-4-yl)penta-2,4-dienylidene)-4,5-dihydro-5-oxopyrazol-1-yl) benzenesulfonate; trisodium 4-(3-ethoxycarbonyl-4-(5-(3-ethoxycarbonyl-5-oxido-1-(4-sulfonatophenyl)pyrazol-4-yl)penta-2,4-dienylidene)-4,5-dihydro-5-oxopyrazol-1-yl)benzenesulfonate	402-660-9	—	Repr. 1B Aquatic Chronic 3	H360-D *** H412	GHS08 Dgr	H360D *** H412			
607-488-00-X	ethyl (2-acetylamino-5-fluoro-4-isothiocyanatophenoxy)acetate	414-210-9	147379-38-2	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-489-00-5	reaction mass of: 2-ethylhexyl linolenate, linoleate and oleate; 2-ethylhexyl epoxyoleate; 2-ethylhexyl diepoxylinoleate; 2-ethylhexyl triepoxylinolenate	414-890-7	71302-79-9	Skin Sens. 1	H317	GHS07 Wng	H317			
607-490-00-0	N-[2-hydroxy-3-(C_{12-16}-alkyloxy)propyl]-N-methyl glycinate	415-060-7	—	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
607-492-00-1	2-(1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy)-2-methyl propyl propanoate	415-490-5	141773-73-1	Aquatic Chronic 2	H411	GHS09	H411			
607-493-00-7	methyl (3a*R*,4*R*,7a*R*)-2-methyl-4-(1*S*,2*R*,3-triacetoxypropyl)-3a,7a-dihydro-4*H*-pyrano[3,4-d]oxazole-6-carboxylate	415-670-3	78850-37-0	Eye Dam. 1	H318	GHS05 Dgr	H318			
607-494-00-2	bis(2-ethylhexyl)octylphosphonate	417-170-0	52894-02-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
607-495-00-8	sodium 4-sulfophenyl-6-((1-oxononyl) amino)hexanoate	417-550-6	168151-92-6	Skin Sens. 1	H317	GHS07 Wng	H317			
607-496-00-3	2,2'-methylenebis(4,6-di-*tert*-butyl-phenyl)-2-ethylhexyl phosphite	418-310-3	126050-54-2	Aquatic Chronic 4	H413	—	H413			
607-497-00-9	cerium oxide isostearate	419-760-3	—	Aquatic Chronic 4	H413	—	H413			
607-498-00-4	(*E*)-3,7-dimethyl-2,6-octadienylhexadecanoate	421-370-3	3681-73-0	Skin Irrit. 2 Aquatic Chronic 4	H315 H413	GHS07 Wng	H315 H413			
607-499-00-X	bis(dimethyl-(2-hydroxyethyl)ammonium) 1,2-ethanediyl-bis(2-hexadecenylsuccinate)	421-660-1	—	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-500-00-3	calcium 2,2,bis[(5-tetrapropylene-2-hydroxy)phenyl]ethanoate	421-670-4	—	Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H315 H400 H410	GHS07 GHS09 Wng	H315 H410			
607-501-00-9	reaction mass of: triphenylthiophosphate and tertiary butylated phenyl derivatives	421-820-9	192268-65-8	Aquatic Chronic 4	H413	—	H413			
607-502-00-4	(*N*-benzyl-N,N,N-tributyl)ammonium 4-dodecylbenzenesulfonate	422-200-0	178277-55-9	Skin Corr. 1B Acute Tox. 4 * Aquatic Chronic 2	H314 H302 H411	GHS05 GHS07 GHS09 Dgr	H314 H302 H411			
607-503-00-X	2,4,6-tri-*n*-propyl-2,4,6-trioxo-1,3,5,2,4,6-trioxatriphosphorinane	422-210-5	68957-94-8	Skin Corr. 1B	H314	GHS05 Dgr	H314			
607-505-00-0	pentasodium 7-(4-(4-(5-amino-4-sulfonato-2-(4-((2-(sulfonato-ethoxy)sulfonyl)phenylazo)phenylamino)-6-chloro-1,3,5-triazin-2-yl)amino-2-ureidophenylazo)naphtalene-1,3,6-trisulfonate	422-930-1		Aquatic Chronic 3	H412	—	H412			
607-506-00-6	reaction mass of: strontium (4-chloro-2-((4,5-dihydro-3-methyl-5-oxo-1-(3-sulfonatophenyl)-1*H*-pyrazol-4-yl)azo)-5-methyl) benzenesulfonate; disodium (4-chloro-2-((4,5-dihydro-3-methyl-5-oxo-1-(3-sulfonatophenyl)-1*H*-pyrazol-4-yl)azo)-5-methyl)benzenesulfonate	422-970-8		Aquatic Chronic 2	H411	GHS09	H411			
607-507-00-1	potassium,sodium 2,4-diamino-3-[4-(2-sulfonatoethoxysulfonyl)phenylazo]-5-[4-(2-sulfonatoethoxysulfonyl)-2-sulfonatophenylazo]-benzenesulfonate	422-980-2	187026-95-5	Eye Dam. 1	H318	GHS05 Dgr	H318			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-508-00-7	disodium 3,3'-[iminobis[sulfonyl-4,1-phenylene-(5-hydroxy-3-methylpyrazole-1,4-diyl)azo-4,1-phenylenesulfonylimino-(4-amino-6-hydroxypyrimidine-2,5-diyl)azo-4,1-phenylenesulfonylimino(4-amino-6-hydroxypyrimidine-2,5-diyl)azo]bis(benzenesulfonate)]	423-110-4	—	Eye Dam. 1	H318	GHS05 Dgr	H318			
607-512-00-9	trisodium 2,4-diamino-3,5-bis-[4-(2-sulfonatoethoxy)sulfonyl)phenylazo]benzenesulfonate	423-970-0	182926-43-8	Aquatic Chronic 3	H412	—	H412			
607-513-00-4	reaction mass of: trisodium 4-benzoylamino-6-(6-ethenesulfonyl-1-sulfato-naphthalen-2-ylazo)-5-hydroxynaphthalene-2,7-disulfonate; 5-(benzoylamino)-4-hydroxy-3-((1-sulfo-6-((2-(sulfooxy)ethyl)sulfonyl)-2-naphthyl)azo) naphthalene-2,7-disulfonic acid sodium salt; 5-(benzoylamino)-4-hydroxy-3-((1-sulfo-6-((2-(sulfooxy)ethyl)sulfonyl)-2-naphthyl)azo) naphthalene-2,7-disulfonic acid	423-200-3	—	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H318 H317 H412	GHS05 GHS07 Dgr	H318 H317 H412			
607-515-00-5	reaction mass of: disodium hexyldiphenyl ether disulphonate; disodium dihexyldiphenyl ether disulphonate	429-650-7	147732-60-3	Eye Irrit. 2 Aquatic Chronic 2	H319 H411	GHS07 GHS09 Wng	H319 H411			
607-516-00-0	*N,N'*-bis(trifluoroacetyl)-*S,S'*-bis-L-homocysteine	429-670-6	105996-54-1	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
607-517-00-6	(*S*)-α-(acetylthio)benzenepropanoic acid	430-300-0	76932-17-7	Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1	H302 H318 H317	GHS05 GHS07 Dgr	H302 H318 H317			
607-526-00-5	cartap (ISO); 1,3-bis(carbamoylthio)-2-(dimethylamino) propane	—	15263-53-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
607-527-00-0	reaction mass of: 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-tridecafluorooctyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heptadecafluorodecyl) dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heneicosafluorododecyl) dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-pentacosafluorotetradecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-heptadecafluorodecyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heptadecafluorodecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-heptadecafluorodecyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heneicosafluorododecyl)dodecanedioate	423-180-6	—	STOT RE 2 *	H373 **	GHS08 Wng	H373 **			
607-696-00-0	pentyl formate	211-340-6	638-49-3	Flam. Liq. 3 Eye Irrit. 2 STOT SE 3	H226 H319 H335	GHS02 GHS07 Dgr	H226 H319 H335			C
607-697-00-6	*tert*-butyl propionate	—	20487-40-5	Flam. Liq. 2	H225	GHS02 Dgr	H225			C
608-001-00-3	acetonitrile; cyanomethane	200-835-2	75-05-8	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2	H225 H332 H312 H302 H319	GHS02 GHS07 Dgr	H225 H332 H312 H302 H319			
608-002-00-9	trichloroacetonitrile	208-885-7	545-06-2	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H331 H311 H301 H411	GHS06 GHS09 Dgr	H331 H311 H301 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-003-00-4	acrylonitrile	203-466-5	107-13-1	Flam. Liq. 2 Carc. 1B Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT SE 3 Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H225 H350 H331 H311 H301 H335 H315 H318 H317 H411	GHS02 GHS06 GHS08 GHS05 GHS09 Dgr	H225 H350 H331 H311 H301 H335 H315 H318 H317 H411		*	D
608-004-00-X	2-hydroxy-2-methylpropionitrile; 2-cyanopropan-2-ol; acetone cyanohydrin	200-909-4	75-86-5	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Aquatic Acute 1 Aquatic Chronic 1	H330 H310 H300 H400 H410	GHS06 GHS09 Dgr	H330 H310 H300 H410			
608-005-00-5	*n*-butyronitrile	203-700-6	109-74-0	Flam. Liq. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H225 H331 H311 H301	GHS02 GHS06 Dgr	H225 H331 H311 H301			
608-006-00-0	bromoxynil (ISO) 3,5-dibromo-4-hydroxybenzonitrile; bromoxynil phenol	216-882-7	1689-84-5	Repr. 2 Acute Tox. 2 * Acute Tox. 3 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H330 H301 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H330 H301 H317 H410		M = 10	

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-007-00-6	ioxynil (ISO) 4-hydroxy-3,5-diiodobenzonitrile	216-881-1	1689-83-4	Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H331 H301 H312 H373 ** H319 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H331 H301 H312 H373 ** H319 H410		M = 10	
608-008-00-1	chloroacetonitrile	203-467-0	107-14-2	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H331 H311 H301 H411	GHS06 GHS09 Dgr	H331 H311 H301 H411			
608-009-00-7	malononitrile	203-703-2	109-77-3	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H410			
608-010-00-2	methacrylonitrile; 2-methyl-2-propene nitrile	204-817-5	126-98-7	Flam. Liq. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Skin Sens. 1	H225 H331 H311 H301 H317	GHS02 GHS06 Dgr	H225 H331 H311 H301 H317		* Skin Sens. 1; H317: C ≥ 0,2 %	D
608-011-00-8	oxalonitrile; cyanogen	207-306-5	460-19-5	Flam. Gas 1 Press. Gas Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H220 H331 H400 H410	GHS02 GHS04 GHS06 GHS09 Dgr	H220 H331 H410			U

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-012-00-3	benzonitrile	202-855-7	100-47-0	Acute Tox. 4 * Acute Tox. 4 *	H312 H302	GHS07 Wng	H312 H302			
608-013-00-9	2-chlorobenzonitrile	212-836-5	873-32-5	Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2	H312 H302 H319	GHS07 Wng	H312 H302 H319			
608-014-00-4	chlorothalonil (ISO); tetrachloroisophthalonitrile	217-588-1	1897-45-6	Carc. 2 Acute Tox. 2 * Eye Dam. 1 STOT SE 3 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H330 H318 H335 H317 H400 H410	GHS06 GHS08 GHS05 GHS09 Dgr	H351 H330 H318 H335 H317 H410		M = 10	
608-015-00-X	dichlobenil (ISO); 2,6-dichlorobenzonitrile	214-787-5	1194-65-6	Acute Tox. 4 * Aquatic Chronic 2	H312 H411	GHS07 GHS09 Wng	H312 H411			
608-016-00-5	1,4-Dicyano-2,3,5,6-tetra-chloro-benzene	401-550-8	1897-41-2	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
608-017-00-0	bromoxynil octanoate (ISO); 2,6-dibromo-4-cyanophenyl octanoate	216-885-3	1689-99-2	Repr. 2 Acute Tox. 3 * Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H331 H302 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H331 H302 H317 H410		M = 10	

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-018-00-6	ioxynil octanoate (ISO); 4-cyano-2,6-diiodophenyl octanoate	223-375-4	3861-47-0	Repr. 2 Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H301 H319 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H301 H319 H317 H410		M = 10	
608-019-00-1	2,2'-dimethyl-2,2'-azodipropiononitrile; ADZN	201-132-3	78-67-1	Self-react. C Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H242 H332 H302 H412	GHS02 GHS07 Dgr	H242 H332 H302 H412			T
608-021-00-2	3-(2-(diaminomethyleneamino)thiazol-4-ylmethylthio)propionitrile	403-710-2	76823-93-3	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			
608-022-00-8	3,7-dimethyloctanenitrile	403-620-3	40188-41-8	Skin Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H315 H317 H411	GHS07 GHS09 Wng	H315 H317 H411			
608-023-00-3	fenbuconazole (ISO); 4-(4-chlorophenyl)-2-phenyl-2-[(1*H*-1,2,4-triazol-1-yl)methyl]butanenitrile	406-140-2	114369-43-6	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
608-024-00-9	2-(4-(*N*-butyl-*N*-phenethylamino)phenyl) ethylene-1,1,2-tricarbonitrile	407-650-8	97460-76-9	Aquatic Chronic 4	H413	—	H413			
608-025-00-4	2-nitro-4,5-bis(benzyloxy)phenylacetonitrile	410-970-0	117568-27-1	Aquatic Chronic 4	H413	—	H413			
608-026-00-X	3-cyano-3,5,5-trimethylcyclohexanone	411-490-4	7027-11-4	Acute Tox. 4 * STOT RE 2 * Skin Sens. 1 Aquatic Chronic 3	H302 H373 ** H317 H412	GHS08 GHS07 Wng	H302 H373 ** H317 H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-027-00-5	reaction mass of: 3-(4-ethylphenyl)-2,2-dimethylpropanenitrile; 3-(2-ethylphenyl)-2,2-dimethylpropanenitrile; 3-(3-ethylphenyl)-2,2-dimethylpropanenitrile	412-660-0	—	Aquatic Chronic 2	H411	GHS09	H411			
608-028-00-0	4-(2-cyano-3-phenylamino acryloyloxymethyl)-cyclohexyl-methyl 2-cyano-3-phenylamino)-acrylate	413-510-7	147374-67-2	STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H373 ** H317 H411	GHS08 GHS09 Wng	H373 ** H317 H411			
608-029-00-6	1,2-dihydro-6-hydroxy-4-methyl-1-[3-(1-methylethoxy)propyl]-2-oxo-3-pyridinecarbonitrile	411-990-2	68612-94-2	Skin Sens. 1	H317	GHS07 Wng	H317			
608-030-00-1	N-acetyl-N-[5-cyano-3-(2-dibutylamino-4-phenylthyazol-5-yl-methylene)-4-methyl-2,6-dioxo-1,2,3,6-tetrahydropyridin-1-yl] benzamide	412-340-0	147741-93-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
608-031-00-7	2-benzyl-2-methyl-3-butenitrile	407-870-4	97384-48-0	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
608-033-00-8	N-butyl-3-(2-chloro-4-nitrophenylhydrazono)-1-cyano-2-methylprop-1-ene-1,3-dicarboximide	407-970-8	75511-91-0	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
608-034-00-3	chlorfenapyr; 4-bromo-2-(4-chlorophenyl)-1-ethoxymethyl-5-trifluoromethylpyrrole-3-carbonitrile	—	122453-73-0	Acute Tox. 3 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H331 H302 H400 H410	GHS06 GHS09 Dgr	H331 H302 H410			
608-035-00-9	(±)-α-[(2-acetyl-5-methylphenyl)-amino]-2,6-dichlorobenzene-aceto-nitrile	419-290-9	—	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			

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608-036-00-4	3-(2-{4-[2-(4-cyanophenyl)vinyl]phenyl} vinyl)benzonitrile	419-060-8	79026-02-1	Aquatic Chronic 4	H413	—	H413			
608-037-00-X	reaction mass of: (*E*)-2,12-tridecadiennitrile; (*E*)-3,12-tridecadiennitrile; (Z)-3,12-tridecadiennitrile	422-190-8		Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
608-038-00-5	2,2,4-trimethyl-4-phenyl-butane-nitrile	422-580-8	75490-39-0	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
608-039-00-0	2-phenylhexanenitrile	423-460-8	3508-98-3	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
608-040-00-6	4,4'-dithiobis(5-amino-1-(2,6-dichloro-4-(trifluoromethyl)phenyl)-1*H*-pyrazole-3-carbonitrile)	423-490-1	130755-46-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
608-041-00-1	4'-((2-butyl-4-oxo-1,3-diazaspiro[4.4]non-1-ene-3-yl)methyl)(1,1'-biphenyl)-2-carbonitrile	423-500-4	138401-24-8	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
608-043-00-2	3-(*cis*-3-hexenyloxy)propanenitril	415-220-6	142653-61-0	Acute Tox. 3 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H331 H302 H400 H410	GHS06 GHS09 Dgr	H331 H302 H410			
608-065-00-2	salts of bromoxynil with the exception of those specified elsewhere in this Annex	—	—	Repr. 2 Acute Tox. 2 * Acute Tox. 3 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H330 H301 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H330 H301 H317 H410		M = 10	A

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
608-066-00-8	salts of ioxynil with the exception of those specified elsewhere in this Annex	—	—	Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H361d *** H331 H301 H312 H373 ** H319 H400 H410	GHS06 GHS08 GHS09 Dgr	H361d *** H331 H301 H312 H373 ** H319 H410		M = 10	A
609-001-00-6	1-nitropropane	203-544-9	108-03-2	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H226 H332 H312 H302	GHS02 GHS07 Wng	H226 H332 H312 H302		*	
609-002-00-1	2-nitropropane	201-209-1	79-46-9	Flam. Liq. 3 Carc. 1B Acute Tox. 4 * Acute Tox. 4 *	H226 H350 H332 H302	GHS02 GHS08 GHS07 Dgr	H226 H350 H332 H302			
609-003-00-7	nitrobenzene	202-716-0	98-95-3	Carc. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Aquatic Chronic 2	H351 H361f *** H331 H311 H301 H372 ** H411	GHS06 GHS08 GHS09 Dgr	H351 H361f *** H331 H311 H301 H372 ** H411			
609-004-00-2	dinitrobenzene; [1] 1,4-dinitrobenzene; [2] 1,3-dinitrobenzene; [3] 1,2-dinitrobenzene [4]	246-673-6 [1] 202-833-7 [2] 202-776-8 [3] 208-431-8 [4]	25154-54-5 [1] 100-25-4 [2] 99-65-0 [3] 528-29-0 [4]	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H330 H310 H300 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H330 H310 H300 H373 ** H410			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-005-00-8	1,3,5-trinitrobenzene	202-752-7	99-35-4	Expl. 1.1 Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H201 H330 H310 H300 H373 ** H400 H410	GHS01 GHS06 GHS08 GHS09 Dgr	H201 H330 H310 H300 H373 ** H410			
609-006-00-3	4-nitrotoluene	202-808-0	99-99-0	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411			
609-007-00-9	2,4-dinitrotoluene; dinitrotoluene, technical grade; [1] dinitrotoluene [2]	204-450-0 [1] 246-836-1 [2]	121-14-2 [1] 25321-14-6 [2]	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H350 H341 H361f *** H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H411			
609-008-00-4	2,4,6-trinitrotoluene; TNT	204-289-6	118-96-7	Expl. 1.1 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H201 H331 H311 H301 H373 ** H411	GHS01 GHS06 GHS08 GHS09 Dgr	H201 H331 H311 H301 H373 ** H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-009-00-X	2,4,6-trinitrophenol; picric acid	201-865-9	88-89-1	Expl. 1.1 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H201 H331 H311 H301	GHS01 GHS06 Dgr	H201 H331 H311 H301			
609-010-00-5	salts of picric acid	—	—	Unst. Expl Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H201 H331 H311 H301	GHS01 GHS06 Dgr	H201 H331 H311 H301			T
609-011-00-0	2,4,6-trinitroanisole	—	606-35-9	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 2	H201 H332 H312 H302 H411	GHS01 GHS07 GHS09 Wng	H201 H332 H312 H302 H411			
609-012-00-6	2,4,6-trinitro-*m*-cresol	210-027-1	602-99-3	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H201 H332 H312 H302	GHS01 GHS07 Wng	H201 H332 H312 H302			
609-013-00-1	2,4,6-trinitro-*m*-xylene	211-187-5	632-92-8	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 *	H201 H332 H312 H302 H373 **	GHS01 GHS08 GHS07 Wng	H201 H332 H312 H302 H373 **			
609-015-00-2	4-nitrophenol; *p*-nitrophenol	202-811-7	100-02-7	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 *	H332 H312 H302 H373 **	GHS08 GHS07 Wng	H332 H312 H302 H373 **			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-016-00-8	dinitrophenol (reaction mass of isomers); [1] 2,4(or 2,6)-dinitrophenol [2]	247-096-2 [1] 275-732-9 [2]	25550-58-7 [1] 71629-74-8 [2]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			
609-018-00-9	2,4,6-trinitroresorcinol; styphnic acid	201-436-6	82-71-3	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H201 H332 H312 H302	GHS01 GHS07 Wng	H201 H332 H312 H302			T
609-019-00-4	lead 2,4,6-trinitro-*m*-phenylene dioxide; lead 2,4,6-trinitroresorcinoxide; lead styphnate	239-290-0	15245-44-0	Unst. Expl Repr. 1A Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H200 H360-Df H332 H302 H373 ** H400 H410	GHS01 GHS08 GHS07 GHS09 Dgr	H200 H360Df H332 H302 H373 ** H410			1
609-019-01-1	lead 2,4,6-trinitro-*m*-phenylene dioxide; lead 2,4,6-trinitroresorcinoxide; lead styphnate (≥ 20 % phlegmatiser)	239-290-0	15245-44-0	Expl. 1.1 Repr. 1A Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H201 H360-Df H332 H302 H373 ** H400 H410	GHS01 GHS08 GHS07 GHS09 Dgr	H201 H360Df H332 H302 H373 ** H410			1

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-020-00-X	DNOC (ISO); 4,6-dinitro-*o*-cresol	208-601-1	534-52-1	Muta. 2 Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H341 H330 H310 H300 H315 H318 H317 H400 H410	GHS06 GHS08 GHS05 GHS07 GHS09 Dgr	H341 H330 H310 H300 H315 H318 H317 H410	EUH044		
609-021-00-5	sodium salt of DNOC; sodium 4,6-dinitro-*o*-cresolate; [1] potassium salt of DNOC; potassium 4,6-dinitro-*o*-cresolate [2]	219-007-7 [1] - [2]	2312-76-7 [1] 5787-96-2 [2]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			
609-022-00-0	ammonium salt of DNOC; ammonium 4,6-dinitro-*o*-tolyl oxide	221-037-0	2980-64-5	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H330 H310 H300 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H330 H310 H300 H373 ** H410			
609-023-00-6	dinocap (ISO)	254-408-0	39300-45-3	Repr. 1B Acute Tox. 4 * STOT RE 2 * Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H332 H373 ** H315 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H360D *** H332 H373 ** H315 H317 H410			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-024-00-1	binapacryl (ISO); 2-*sec*-butyl-4,6-dinitrophenyl-3-methylcrotonate	207-612-9	485-31-4	Repr. 1B Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H312 H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H360D *** H312 H302 H410			
609-025-00-7	dinoseb(ISO); 6-*sec*-butyl-2,4-dinitrophenol	201-861-7	88-85-7	Repr. 1B Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H360-Df H311 H301 H319 H400 H410	GHS06 GHS08 GHS09 Dgr	H360Df H311 H301 H319 H410	EUH044		
609-026-00-2	salts and esters of dinoseb, with the exception of those specified elsewhere in this Annex	—	—	Repr. 1B Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H360-Df H311 H301 H319 H400 H410	GHS06 GHS08 GHS09 Dgr	H360Df H311 H301 H319 H410	EUH044		A
609-027-00-8	dinocton; reaction mass of isomers: methyl 2-octyl-4,6-dinitrophenyl carbonate, methyl 4-octyl-2,6-dinitrophenyl carbonate	—	63919-26-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
609-028-00-3	dinex (ISO); 2-cyclohexyl-4,6-dinitrophenol	205-042-5	131-89-5	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-029-00-9	salts and esters of dinex	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H410			A
609-030-00-4	dinoterb (ISO); 2-*tert*-butyl-4,6-dinitrophenol	215-813-8	1420-07-1	Repr. 1B Acute Tox. 2 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H300 H311 H400 H410	GHS06 GHS08 GHS09 Dgr	H360D *** H300 H311 H410	EUH044		
609-031-00-X	salts and esters of dinoterb	—	—	Repr. 1B Acute Tox. 2 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H300 H311 H400 H410	GHS06 GHS08 GHS09 Dgr	H360D *** H300 H311 H410			A
609-032-00-5	bromofenoxim (ISO); 3,5-dibromo-4-hydroxybenzaldehyde-O-(2,4-dinitrophenyl)-oxime	236-129-6	13181-17-4	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
609-033-00-0	dinosam (ISO); 2-(1-methylbutyl)-4,6-dinitrophenol	—	4097-36-3	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-034-00-6	salts and esters of dinosam	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H410			A
609-035-00-1	nitroethane	201-188-9	79-24-3	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 *	H226 H332 H302	GHS02 GHS07 Wng	H226 H332 H302		*	
609-036-00-7	nitromethane	200-876-6	75-52-5	Flam. Liq. 3 Acute Tox. 4 *	H226 H302	GHS02 GHS07 Wng	H226 H302		*	
609-037-00-2	5-nitroacenaphthene	210-025-0	602-87-9	Carc. 1B	H350	GHS08 Dgr	H350			
609-038-00-8	2-nitronaphthalene	209-474-5	581-89-5	Carc. 1B Aquatic Chronic 2	H350 H411	GHS08 GHS09 Dgr	H350 H411			
609-039-00-3	4-nitrobiphenyl	202-204-7	92-93-3	Carc. 1B Aquatic Chronic 2	H350 H411	GHS08 GHS09 Dgr	H350 H411			
609-040-00-9	nitrofen (ISO); 2,4-dichlorophenyl 4-nitrophenyl ether	217-406-0	1836-75-5	Carc. 1B Repr. 1B Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H350 H360-D *** H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H360D *** H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-041-00-4	2,4-dinitrophenol	200-087-7	51-28-5	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1	H331 H311 H301 H373 ** H400	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H400			
609-042-00-X	pendimethalin (ISO); N-(1-ethylpropyl)-2,6-dinitro-3,4-xylidine	254-938-2	40487-42-1	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
609-043-00-5	quintozene (ISO); pentachloronitrobenzene	201-435-0	82-68-8	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
609-044-00-0	tecnazene (ISO); 1,2,4,5-tetrachloro-3-nitrobenzene	204-178-2	117-18-0	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
609-045-00-6	reaction mass of: 4,6-dinitro-2-(3-octyl) phenyl methyl carbonate and 4,6-dinitro-2-(4-octyl)phenyl methyl carbonate; dinocton-6	—	8069-76-9	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
609-046-00-1	trifluralin (ISO) (containing < 0.5 ppm NPDA); α,α,α-trifluoro-2,6-dinitro-*N*,*N*-dipropyl-*p*-toluidine (containing < 0.5 ppm NPDA); 2,6-dinitro-*N*,*N*-dipropyl-4-trifluoromethylaniline (containing < 0.5 ppm NPDA); *N*,*N*-dipropyl-2,6-dinitro-4-trifluoromethylaniline (containing < 0.5 ppm NPDA)	216-428-8	1582-09-8	Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H319 H317 H400 H410	GHS07 GHS09 Wng	H319 H317 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-047-00-7	2-nitroanisole	202-052-1	91-23-6	Carc. 1B Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			
609-048-00-2	sodium 3-nitrobenzenesulphonate	204-857-3	127-68-4	Eye Irrit. 2 Skin Sens. 1	H319 H317	GHS07 Wng	H319 H317			
609-049-00-8	2,6-dinitrotoluene	210-106-0	606-20-2	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 3	H350 H341 H361f *** H331 H311 H301 H373 ** H412	GHS06 GHS08 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H412			
609-050-00-3	2,3-dinitrotoluene	210-013-5	602-01-7	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H350 H341 H361f *** H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-051-00-9	3,4-dinitrotoluene	210-222-1	610-39-9	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H350 H341 H361f *** H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H411			
609-052-00-4	3,5-dinitrotoluene	210-566-2	618-85-9	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 3	H350 H341 H361f *** H331 H311 H301 H373 ** H412	GHS06 GHS08 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H412			
609-053-00-X	hydrazine-trinitromethane	414-850-9	—	Expl. 1.1 **** Self-react. A Carc. 1B Acute Tox. 3 * Acute Tox. 3 * Skin Sens. 1	H201 H240 H350 H331 H301 H317	GHS01 GHS06 GHS08 Dgr	H201 H240 H350 H331 H301 H317			
609-054-00-5	2,3-dinitrophenol; [1] 2,5-dinitrophenol; [2] 2,6-dinitrophenol; [3] 3,4-dinitrophenol; [4] salts of dinitrophenol [5]	200-628-7 [1] 206-348-1 [2] 209-357-9 [3] 209-415-3 [4] - [5]	66-56-8 [1] 329-71-5 [2] 573-56-8 [3] 577-71-9 [4] - [5]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-055-00-0	2,5-dinitrotoluene	210-581-4	619-15-8	Carc. 1B Muta. 2 Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H350 H341 H361f *** H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H350 H341 H361f *** H331 H311 H301 H373 ** H411			
609-056-00-6	2,2-dibromo-2-nitroethanol	412-380-9	69094-18-4	Expl. 1.1 Carc. 2 Acute Tox. 4 * STOT RE 2 * Skin Corr. 1A Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H201 H351 H302 H373 ** H314 H317 H400 H410	GHS01 GHS08 GHS05 GHS07 GHS09 Dgr	H201 H351 H302 H373 ** H314 H317 H410		* STOT SE 3; H335: C ≥ 1 %	T
609-057-00-1	3-chloro-2,4-difluoronitrobenzene	411-980-8	3847-58-3	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H314 H317 H410			
609-058-00-7	2-nitro-2-phenyl-1,3-propanediol	410-360-4	5428-02-4	STOT RE 1 Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H372 ** H312 H302 H317 H411	GHS08 GHS07 GHS09 Dgr	H372 ** H312 H302 H317 H411	EUH070		

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-059-00-2	2-chloro-6-(ethylamino)-4-nitrophenol	411-440-1	131657-78-8	Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H302 H317 H411	GHS07 GHS09 Wng	H302 H317 H411			
609-060-00-8	4-[(3-hydroxypropyl)amino]-3-nitrophenol	406-305-9	92952-81-3	Skin Irrit. 2 Aquatic Chronic 2	H315 H411	GHS07 GHS09 Wng	H315 H411			
609-061-00-3	(*E*,*Z*)-4-chlorophenyl(cyclopropyl)ketone *O*-(4-nitrophenylmethyl)oxime	406-100-4	94097-88-8	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
609-062-00-9	2-bromo-2-nitropropanol	407-030-7	24403-04-1	Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H311 H302 H373 ** H314 H317 H400 H410	GHS06 GHS08 GHS05 GHS09 Dgr	H311 H302 H373 ** H314 H317 H410			
609-063-00-4	2-[(4-chloro-2-nitrophenyl)amino]ethanol	413-280-8	59320-13-7	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
609-064-00-X	mesotrione(ISO); 2-[4-(methylsulfonyl)-2-nitrobenzoyl]-1,3-cyclohexanedione	—	104206-82-8	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
609-065-00-5	2-nitrotoluene	201-853-3	88-72-2	Carc. 1B Muta. 1B Repr. 2 Acute Tox. 4 * Aquatic Chronic 2	H350 H340 H361f *** H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H340 H361f *** H302 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
609-066-00-0	lithium sodium 3-amino-10-{4-(10-amino-6,13-dichloro-4,11-disulfonatobenzo[5,6][1,4]oxazino[2,3-b]phenoxazine-3-ylamino)-6-[methyl(2-sulfonato-ethyl)amino]-1,3,5-triazin-2-ylamino}-6,13-dichlorobenzo[5,6][1,4]oxazino[2,3-b]phenoxazine-4,11-disulfonate	418-870-9	154212-58-5	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * STOT SE 2 **	H332 H312 H302 H371 **	GHS08 GHS07 Dgr	H332 H312 H302 H371 **			
609-067-00-6	sodium and potassium 4-(3-aminopropylamino)-2,6-bis[3-(4-methoxy-2-sulfophenylazo)-4-hydroxy-2-sulfo-7-naphthylamino]-1,3,5-triazine	416-280-6	156769-97-0	Skin Sens. 1	H317	GHS07 Wng	H317			
609-068-00-1	musk xylene; 5-*tert*-butyl-2,4,6-trinitro-*m*-xylene	201-329-4	81-15-2	Expl. 1.1 Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H201 H351 H400 H410	GHS01 GHS08 GHS09 Wng	H201 H351 H410			T
609-070-00-2	1,4-dichloro-2-(1,1,2,3,3,3-hexafluoropropoxy)-5-nitrobenzene	415-580-4	130841-23-5	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
609-071-00-8	reaction mass of: 2-methylsulfanyl-4,6-bis-(2-hydroxy-4-methoxy-phenyl)-1,3,5-triazine; 2-(4,6-bis-methylsulfanyl-1,3,5-triazin-2-yl)-5-methoxy-phenol	423-520-3	156137-33-6	Skin Sens. 1	H317	GHS07 Wng	H317			
610-001-00-3	trichloronitromethane; chloropicrin	200-930-9	76-06-2	Acute Tox. 2 * Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2	H330 H302 H319 H335 H315	GHS06 Dgr	H330 H302 H319 H335 H315			
610-002-00-9	1,1-dichloro-1-nitroethane	209-854-0	594-72-9	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H331 H311 H301	GHS06 Dgr	H331 H311 H301			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
610-003-00-4	chlorodinitrobenzene	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			C
610-004-00-X	2-chloro-1,3,5-trinitrobenzene	201-864-3	88-88-0	Expl. 1.1 Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Aquatic Acute 1 Aquatic Chronic 1	H201 H330 H310 H300 H400 H410	GHS01 GHS06 GHS09 Dgr	H201 H330 H310 H300 H410			
610-005-00-5	1-chloro-4-nitrobenzene	202-809-6	100-00-5	Carc. 2 Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H351 H341 H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H351 H341 H331 H311 H301 H373 ** H411			
610-006-00-0	chloronitroanilines with the exception of those specified elsewhere in this Annex	—	—	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Chronic 2	H330 H310 H300 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H330 H310 H300 H373 ** H411			A C
610-007-00-6	1-chloro-1-nitropropane	209-990-0	600-25-9	Acute Tox. 4 * Acute Tox. 4 *	H332 H302	GHS07 Wng	H332 H302		*	
610-008-00-1	2,6-dichloro-4-nitroanisole	403-350-6	17742-69-7	Acute Tox. 3 * Aquatic Chronic 2	H301 H411	GHS06 GHS09 Dgr	H301 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
610-009-00-7	2-chloro-4-nitroaniline	204-502-2	121-87-9	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
610-010-00-2	2-bromo-1-(2-furyl)-2-nitroethylene	406-110-9	35950-52-8	Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H373 ** H314 H317 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H302 H373 ** H314 H317 H410			
611-001-00-6	azobenzene	203-102-5	103-33-3	Carc. 1B Muta. 2 Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H350 H341 H332 H302 H373 ** H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H341 H332 H302 H373 ** H410			
611-002-00-1	azoxybenzene	207-802-1	495-48-7	Acute Tox. 4 * Acute Tox. 4 *	H332 H302	GHS07 Wng	H332 H302			
611-003-00-7	fenaminosulf (ISO); sodium 4-dimethylaminobenzenediazosulphonate	205-419-4	140-56-7	Acute Tox. 3 * Acute Tox. 4 * Aquatic Chronic 3	H301 H312 H412	GHS06 Dgr	H301 H312 H412			
611-004-00-2	methyl-ONN-azoxymethyl acetate; methyl azoxy methyl acetate	209-765-7	592-62-1	Carc. 1B Repr. 1B	H350 H360D ***	GHS08 Dgr	H350 H360D ***			
611-005-00-8	disodium {5-[(4'-((2,6-hydroxy-3-((2-hydroxy-5-sulphophenyl)azo)phenyl)azo)(1,1'-biphenyl)-4-yl)azo]salicylato(4-)}cuprate(2-); CI Direct Brown 95	240-221-1	16071-86-6	Carc. 1B	H350	GHS08 Dgr	H350			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-006-00-3	4-*o*-tolylazo-*o*-toluidine; 4-amino-2',3-dimethylazobenzene; fast garnet GBC base; AAT; *o*-aminoazotoluene	202-591-2	97-56-3	Carc. 1B Skin Sens. 1	H350 H317	GHS08 Dgr	H350 H317			
611-007-00-9	tricyclazole (ISO); 5-methyl-1,2,4-triazolo(3,4-b)benzo-1,3-thiazole;	255-559-5	41814-78-2	Acute Tox. 4 *	H302	GHS07 Wng	H302			
611-008-00-4	4-aminoazobenzene; 4-phenylazoaniline	200-453-6	60-09-3	Carc. 1B Aquatic Acute 1 Aquatic Chronic 1	H350 H400 H410	GHS08 GHS09 Dgr	H350 H410			
611-009-00-X	sodium (1-(5-(4-(4-anilino-3-sulphophenylazo)-2-methyl-5-methylsulphonamidophenylazo)-4-hydroxy-2-oxido-3-(phenylazo) phenylazo)-5-nitro-4-sulphonato-2-naphtholato)iron(II)	401-220-3	—	Acute Tox. 4 * Aquatic Chronic 3	H332 H412	GHS07 Wng	H332 H412			
611-010-00-5	2'-(2-cyano-4,6-dinitrophenylazo)-5'-(*N*,*N*-dipropylamino)propionanilide	403-010-7	106359-94-8	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
611-011-00-0	*N*,*N*,*N*',*N*'-tetramethyl-3,3'-(propylenebis (iminocarbonyl-4,1-phenylenazo(1,6-dihydro-2-hydroxy-4-methyl-6-oxopyridine-3,1-diyl)))di(propylammonium) dilactate	403-340-1	—	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dg	H318 H411			
611-012-00-6	reaction mass of 2,2-iminodiethanol 6-methyl-2-(4-(2,4,6-triaminopyrimidin-5-ylazo)phenyl)benzothiazole-7-sulfonate and 2-methylaminoethanol 6-methyl-2-(4-(2,4,6-triaminopyrimidin-5-ylazo)phenyl) benzothiazole-7-sulfonate and *N*,*N*-diethylpropane-1,3-diamine 6-methyl-2-(4-(2,4,6-triaminopyrimidin-5-ylazo)phenyl)benzothiazole-7-sulfonate	403-410-1	114565-65-0	Skin Sens. 1	H317	GHS07 Wng	H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-013-00-1	trilithium-1-hydroxy-7-(3-sulfonatoanilino)-2-(3-methyl-4-(2-methoxy-4-(3-sulfonatophenylazo)phenylazo)phenylazo)naphthalene-3-sulfonate	403-650-7	117409-78-6	Expl. 1.3 **** Aquatic Chronic 2	H203 H411	GHS01 GHS09 Dgr	H203 H411			
611-014-00-7	(tetrasodium 1-(4-(3-acetamido-4-(4'-nitro-2,2'-disulfonatostilben-4-ylazo)anilino)-6-(2,5-disulfonatoanilino)-1,3,5-triazin-2-yl)-3-carboxypyridinium) hydroxide	404-250-5	115099-55-3	Skin Sens. 1	H317	GHS07 Wng	H317			
611-015-00-2	tetrasodium 4-amino-5-hydroxy-6-(4-(2-(2-(sulfonatooxy)ethylsulfonyl)ethylcarbamoyl)phenylazo)-3-(4-(2-(sulfonatooxy)ethylsulfonyl)phenylazo)naphthalene-2,7-disulfonate	404-320-5	116889-78-2	Skin Sens. 1	H317	GHS07 Wng	H317			
611-016-00-8	reaction mass of 1,1'-((dihydroxyphenylene) bis(azo-3,1-phenylenazo(1-(3-dimethylaminopropyl)-1,2-dihydro-6-hydroxy-4-methyl-2-oxopyridine-5,3-diyl)))dipyridinium dichloride dihydrochloride, mixed isomers and 1-(1-(3-dimethylaminopropyl)-5-(3-((4-(1-(3-dimethylaminopropyl)-1,6-dihydro-2-hydroxy-4-methyl-6-oxo-5-pyridinio-3-pyridylazo)phenylazo)-2,4(or2,6 or3,5)-dihydroxyphenylazo)phenylazo)-1,2-dihydro-6-hydroxy-4-methyl-2-oxo-3-pyridyl)pyridinium dichloride	404-540-1	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-017-00-3	2-(4-(diethylaminopropylcarbamoyl)phenylazo)-3-oxo-N-(2,3-dihydro-2-oxobenzimidazol-5-yl)butyramide	404-910-2	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-018-00-9	tetraammonium 5-(4-(7-amino-1-hydroxy-3-sulfonato-2-naphthylazo)-6-sulfonato-1-naphthylazo)isophthalate	405-130-5	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-019-00-4	tetralithium 6-amino-4-hydroxy-3-(7-sulfonato-4-(4-sulfonatophenylazo)-1-naphthylazo)naphthalene-2,7-disulfonate	405-150-4	106028-58-4	Skin Sens. 1	H317	GHS07 Wng	H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-020-00-X	tetrakis(tetramethylammonium) 6-amino-4-hydroxy-3-(7-sulfonato-4-(4-sulfonatophenylazo)-1-naphthylazo)naphthalene-2,7-disulfonate	405-170-3	116340-05-7	Acute Tox. 3 * Skin Sens. 1 Aquatic Chronic 3	H301 H317 H412	GHS06 Dgr	H301 H317 H412			
611-021-00-5	2-(4-(4-cyano-3-methylisothiazol-5-ylazo)-N-ethyl-3-methylanilino)ethyl acetate	405-480-9	—	Acute Tox. 4 * STOT RE 2 * Skin Irrit. 2 Aquatic Chronic 4	H302 H373 ** H315 H413	GHS08 GHS07 Wng	H302 H373 ** H315 H413			
611-022-00-0	4-dimethylaminobenzenediazonium 3-carboxy-4-hydroxybenzenesulfonate	404-980-4	—	Self-react. C Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H242 H331 H301 H312 H373 ** H318 H317 H400 H410	GHS02 GHS06 GHS08 GHS05 GHS09 Dgr	H242 H331 H301 H312 H373 ** H318 H317 H410			T
611-023-00-6	disodium 7-(4,6-dichloro-1,3,5-triazin-2-ylamino)-4-hydroxy-3-(4-(2-(sulfonatooxy) ethylsulfonyl)phenylazo) naphthalene-2-sulfonate	404-600-7	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-024-00-1	Benzidine based azo dyes; 4,4'-diarylazobiphenyl dyes, with the exception of those specified elsewhere in this Annex	—	—	Carc. 1B	H350	GHS08 Dgr	H350			A
611-025-00-7	disodium 4-amino-3-[[4'-[(2,4-diaminophenyl)azo][1,1'-biphenyl]-4-yl]azo]-5-hydroxy-6-(phenylazo)naphtalene-2,7-disulphonate; C.I. Direct Black 38	217-710-3	1937-37-7	Carc. 1B Repr. 2	H350 H361d ***	GHS08 Dgr	H350 H361d ***			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-026-00-2	tetrasodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis (azo)]bis[5-amino-4-hydroxynaphthalene-2,7-disulphonate]; C.I. Direct Blue 6	220-012-1	2602-46-2	Carc. 1B Repr. 2	H350 H361d ***	GHS08 Dgr	H350 H361d ***			
611-027-00-8	disodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis (azo)]bis(4-aminonaphthalene-1-sulphonate); C.I. Direct Red 28	209-358-4	573-58-0	Carc. 1B Repr. 2	H350 H361d ***	GHS08 Dgr	H350 H361d ***			
611-028-00-3	C,C'-azodi(formamide)	204-650-8	123-77-3	Resp. Sens. 1	H334	GHS08 Dgr	H334			G
611-029-00-9	*o*-dianisidine based azo dyes; 4,4'-diarylazo-3,3'-dimethoxybiphenyl dyes with the exception of those mentioned elsewhere in this Annex	—	—	Carc. 1B	H350	GHS08 Dgr	H350			A H
611-030-00-4	*o*-tolidine based dyes; 4,4'-diarylazo-3,3'-dimethylbiphenyl dyes, with the exception of those mentioned elsewhere in this Annex	—	—	Carc. 1B	H350	GHS08 Dgr	H350			A H
611-031-00-X	4,4'-(4-iminocyclohexa-2,5-dienylidene-methylene)dianiline hydrochloride; C.I. Basic Red 9	209-321-2	569-61-9	Carc. 1B	H350	GHS08 Dgr	H350			
611-032-00-5	1,4,5,8-tetraaminoanthraquinone; C.I. Disperse Blue 1	219-603-7	2475-45-8	Carc. 1B Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1	H350 H315 H318 H317	GHS08 GHS05 GHS07 Dgr	H350 H315 H318 H317			
611-033-00-0	hexasodium [4,4"-azoxybis(2,2'-disulfonatostilbene-4,4'-diylazo)]-bis[5'-sulfonatobenzene-2,2'- diolato-O(2),O(2),N(1)]-copper(II)	400-020-3	82027-60-9	Aquatic Chronic 2	H411	GHS09	H411			
611-034-00-6	*N*-(5-(bis(2-methoxyethyl)amino)-2-((5-nitro-2,1-benzisothiazol-3-yl)azo)phenylacetamide	402-430-8	105076-77-5	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-035-00-1	tetralithium 6-amino-4-hydroxy-3-[7-sulfonato-4-(5-sulfonato-2-naphthylazo)-1-naphthylazo]naphthalene-2,7-disulfonate	403-660-1	107246-80-0	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-036-00-7	2-(4-(5,6(or 6,7)-dichloro-1,3-benzothiazol-2-ylazo)-*N*-methyl-*m*-toluidino)ethyl acetate	405-440-0	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-037-00-2	3(or 5)-(4-(*N*-benzyl-*N*-ethylamino)-2-methylphenylazo)-1,4-dimethyl-1,2,4-triazolium methylsulphate	406-055-0	124584-00-5	Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H302 H318 H317 H411			
611-038-00-8	trisodium 1-hydroxynaphthalene-2-azo-4'(5',5"-dimethylbiphenyl)-4"-azo(4"-phenylsulfonyloxybenzene)- 2',2",4-trisulfonate	406-820-9	—	Eye Irrit. 2	H319	GHS07 Wng	H319			
611-039-00-3	7-[((4,6-dichloro-1,3,5-triazin-2-yl)amino)-4-hydroxy-3-(4-((2-sulfoxy)ethyl)sulfonyl)phenylazo]naphthalene-2-sulfonic acid	407-050-6	117715-57-8	Skin Sens. 1	H317	GHS07 Wng	H317			
611-040-00-9	3-(5-acetylamino-4-(4-[4,6-bis(3-diethylaminopropylamino)-1,3,5-triazin-2-ylamino]phenylazo)-2-(2-methoxyethoxy)phenylazo)-6-amino-4-hydroxy-2-naphthalenesulfonic acid	407-670-7	115099-58-6	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
611-041-00-4	2-[[4[[4,6-bis[[3-(diethylamino)propyl]amino]-1,3,5-triazine-2-yl]amino]phenyl]azo]-*N*-(2,3-dihydro-2-oxo-1*H*-benzimidazol-5-yl)-3-oxobutanamide	407-680-1	98809-11-1	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			
611-042-00-X	trisodium 5-amino-3-[5-(2-bromoacryloylamino)-2-sulfonatophenylazo]-4-hydroxy-6-(4-vinylsulfonylphenylazo)naphthalene-2,7-disulfonate	411-770-6	136213-71-3	Aquatic Chronic 3	H412	—	H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-043-00-5	reaction mass of: trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6-[2-amino-4-(or 6)-hydroxy-(or 4-amino-2-hydroxy)phenylazo]-6''-(1-carbaniloyl-2-hydroxyprop-1-enylazo)-5',5'''-disulfamoyl-3,3''-disulfonatobis (naphthalene-2,1'-azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate; trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6,6''-bis (1-carbaniloyl-2-hydroxyprop-1-enylazo)-5',5'''-disulfamoyl-3,3''-disulfonatobis (naphthalene-2,1'azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate; trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6,6''-bis [2-amino-4-(or 6)-hydroxy-(or 4-amino-2-hydroxy)phenylazo]5',5'''-disulfamoyl-3,3''-disulfonatobis(naphthalene-2,1'azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate (2:1:1)	402-850-1	—	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			

611-044-00-0	reaction mass of: *tert*-alkyl(C_{12}-C_{14})ammonium bis[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium bis[1-[(2-hydroxy-4-nitrophenyl)azo]-2-naphthalenolato(2-)]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium bis[1-[[5-(1,1-dimethylpropyl)-2-hydroxy-3-nitrophenyl]azo]-2-naphthalenolato(2-)]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium [[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]-[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium [[1-[[5-(1,1-dimethylpropyl)-2-hydroxy-3-nitrophenyl]azo]-2-naphthalenolato(2-)]-[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium ((1-(4(or 5)-nitro-2-oxidophenylazo)-2-naphtholato)(1-(3-nitro-2-oxido-5-pentylphenylazo)-2-naphtholato))chromate(1-)
		403-720-7	117527-94-3	Aquatic Chronic 2	H411	GHS09	H411			

611-045-00-6	2-[4-[*N*-(4-acetoxybutyl)-*N*-ethyl]amino-2-methylphenylazo]-3-acetyl-5-nitrothiophene	404-830-8	—	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-046-00-1	4,4'-diamino-2-methylazobenzene	407-590-2	43151-99-1	Acute Tox. 3 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H301 H373 ** H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H301 H373 ** H317 H410			
611-047-00-7	reaction mass of: 2-[[4-[*N*-ethyl-*N*-(2-acetoxyethyl)amino]phenyl]azo]-5,6-dichlorobenzothiazole; 2-[[4-[*N*-ethyl-*N*-(2-acetoxyethyl)amino]phenyl]azo]-6,7-dichlorobenzothiazole (1:1)	407-890-3	111381-11-4	Aquatic Chronic 4	H413	—	H413			
611-048-00-2	reaction mass of: 2-[[4-[bis(2-acetoxyethyl)amino]phenyl]azo]-5,6-dichlorobenzothiazole; 2-[[4-[bis(2-acetoxyethyl)amino]phenyl]azo]-6,7-dichlorobenzothiazole (1:1)	407-900-6	111381-12-5	Aquatic Chronic 4	H413	—	H413			
611-049-00-8	reaction mass of 7-[4-(3-diethylaminopropylamino)-6-(3-diethylammoniopropylamino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(4-phenylazophenylazo)-naphthalene-2-sulfonate, acetic acid, lactic acid (2:1:1)	408-000-6	118658-98-3	STOT RE 2 * Skin Sens. 1 Aquatic Chronic 3	H373 ** H317 H412	GHS08 Wng	H373 ** H317 H412			
611-051-00-9	2-(4-(*N*-ethyl-*N*-(2-hydroxy)ethyl)amino-2-methylphenyl)azo-6-methoxy-3-methylbenzothiazolium chloride	411-110-7	136213-74-6	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
611-052-00-4	monosodium aqua-[5-[[2,4-dihydroxy-5-[(2-hydroxy-3,5-dinitrophenyl)azo]phenyl]azo]-2-naphthalensulfonate], iron complex	400-720-9	—	Aquatic Chronic 3	H412	—	H412			
611-053-00-X	2,2'-azobis[2-methylpropionamidine] dihydrochloride	221-070-0	2997-92-4	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			
611-055-00-0	C.I. Disperse Yellow 3; *N*-[4-[(2-hydroxy-5-methylphenyl)azo]phenyl]acetamide	220-600-8	2832-40-8	Carc. 2 Skin Sens. 1	H351 H317	GHS08 GHS07 Wng	H351 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-056-00-6	C.I. Solvent Yellow 14; 1-phenylazo-2-naphthol	212-668-2	842-07-9	Carc. 2 Muta. 2 Skin Sens. 1 Aquatic Chronic 4	H351 H341 H317 H413	GHS08 GHS07 Wng	H351 H341 H317 H413			
611-057-00-1	6-hydroxy-1-(3-isopropoxypropyl)-4-methyl-2-oxo-5-[4-(phenylazo)phenylazo]-1,2-dihydro-3-pyridinecarbonitrile	400-340-3	85136-74-9	Carc. 1B Aquatic Chronic 4	H350 H413	GHS08 Wng	H350 H413			
611-058-00-7	(6-(4-hydroxy-3-(2-methoxyphenylazo)-2-sulfonato-7-naphthylamino)-1,3,5-triazin-2,4-diyl)bis[(amino-1-methylethyl)ammonium] formate	402-060-7	108225-03-2	Carc. 1B Eye Dam. 1 Aquatic Chronic 2	H350 H318 H411	GHS08 GHS05 GHS09 Dgr	H350 H318 H411			
611-059-00-2	octasodium 2-(6-(4-chloro-6-(3-(N-methyl-N-(4-chloro-6-(3,5-disulfonato-2-naphthylazo)-1-hydroxy-6-naphthylamino)-1,3,5-triazin-2-yl)aminomethyl)phenylamino)-1,3,5-triazin-2-ylamino)-3,5-disulfonato-1-hydroxy-2-naphthylazo)naphthalene-1,5-disulfonate	412-960-1	148878-21-1	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H318 H317 H412	GHS05 GHS07 Dgr	H318 H317 H412			
611-060-00-8	reaction mass of: sodium 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonatonaphthalen-2-ylazo]-isophthalate; ammonium 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonatonaphthalen-2-ylazo]-isophthalate; 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonaphthalen-2-ylazo]-isophthalic acid	413-180-4	187285-15-0	Eye Dam. 1	H318	GHS05 Dgr	H318			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-061-00-3	disodium 5-[5-[4-(5-chloro-2,6-difluoropyrimidin-4-ylamino)benzamido]-2-sulfonatophenylazo]-1-ethyl-6-hydroxy-4-methyl-2-oxo-3-pyridylmethylsulfonate	412-530-3	—	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
611-062-00-9	octasodium 2-(8-(4-chloro-6-(3-((4-chloro-6-(3,6-disulfonato-2-(1,5-disulfonatonaphthalen-2-ylazo)-1-hydroxynaphthalen-8-ylamino)-1,3,5-triazin-2-yl)aminomethyl)phenylamino)-1,3,5-triazin-2-ylamino)-3,6-disulfonato-1-hydroxynaphthalen-2-ylazo)naphthalene-1,5-disulfonate	413-550-5	—	Skin Irrit. 2 Eye Dam. 1	H315 H318	GHS05 Dgr	H315 H318			
611-063-00-4	trisodium [4'-(8-acetylamino-3,6-disulfonato-2-naphthylazo)-4"-(6-benzoylamino-3-sulfonato-2-naphthylazo)-biphenyl-1,3',3",1'"-tetraolato-*O*,*O*',*O*",*O*'"]copper(II)	413-590-3	164058-22-4	Carc. 1B	H350	GHS08 Dgr	H350			
611-064-00-X	4-(3,4-dichlorophenylazo)-2,6-di-*sec*-butylphenol	410-600-8	124719-26-2	STOT RE 2 * Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H315 H400 H410	GHS08 GHS07 GHS09 Wng	H373 ** H315 H410			
611-065-00-5	4-(4-nitrophenylazo)-2,6-di-*sec*-butyl-phenol	410-610-2	111850-24-9	STOT RE 2 * Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H319 H315 H317 H400 H410	GHS08 GHS07 GHS09 Wng	H373 ** H319 H315 H317 H410			
611-066-00-0	tetrasodium 5-[4-chloro-6-(*N*-ethyl-anilino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(1,5-disulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate	411-540-5	130201-57-9	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-067-00-6	reaction mass of: bis(tris(2-(2-hydroxy(1-methyl)ethoxy)ethyl)ammonium) 7-anilino-4-hydroxy-3-(2-methoxy-5-methyl-4-(4-sulfonatophenylazo)phenylazo)naphthalene-2-sulfonate; bis(tris(2-(2-hydroxy(2-methyl)ethoxy)ethyl) ammonium) 7-anilino-4-hydroxy-3-(2-methoxy-5-methyl-4-(4-sulfonatophenylazo)phenylazo)naphthalene-2-sulfonate	406-910-8	—	Acute Tox. 4 * Eye Dam. 1 Aquatic Chronic 3	H302 H318 H412	GHS05 GHS07 Dgr	H302 H318 H412			
611-068-00-1	tetrasodium 4-amino-3,6-bis(5-[4-chloro-6-(2-hydroxyethylamino)-1,3,5-triazin-2-ylamino]-2-sulfonatophenylazo)-5-hydroxynaphthalene-2,7-disulfonate	400-690-7	85665-98-1	Aquatic Chronic 2	H411	GHS09	H411			
611-069-00-7	*N*,*N*-di-[poly(oxyethylene)-co-poly(oxypropylene)]-4-[(3,5-dicyano-4-methyl-2-thienyl)azo)]-3-methylaniline	413-380-1	—	Aquatic Chronic 2	H411	GHS09	H411			
611-070-00-2	reaction mass of: disodium (6-(4-anisidino)-3-sulfonato-2-(3,5-dinitro-2-oxidophenylazo)-1-naphtholato)(1-(5-chloro-2-oxidophenylazo)-2-naphtholato)chromate(1-); trisodium bis(5-(4-anisidino)-3-sulfonato-2-(3,5-dinitro-2-oxidophenylazo)-1-naphtholato)chromate(1-)	405-665-4	—	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
611-071-00-8	tris(tetramethylammonium) 5-hydroxy-1-(4-sulphonatophenyl)-4-(4-sulphonatophenylazo)pyrazole-3-carboxylate	406-073-9	131013-81-5	Acute Tox. 3 * Aquatic Chronic 3	H301 H412	GHS06 Dgr	H301 H412			
611-072-00-3	2,4-bis[2,2'-[2-(*N*,*N*-dimethylamino)ethyloxycarbonyl]phenylazo]-1,3-dihydroxybenzene, dihydrochloride	407-010-8	118208-02-9	Acute Tox. 4 * Eye Dam. 1 Aquatic Chronic 2	H302 H318 H411	GHS05 GHS07 GHS09 Dgr	H302 H318 H411			
611-073-00-9	dimethyl 3,3'-(*N*-(4-(4-bromo-2,6-dicyanophenylazo)-3-hydroxyphenyl)imino)dipropionate	407-310-9	122630-55-1	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-074-00-4	reaction mass of: sodium/potassium (3-(4-(5-(5-chloro-2,6-difluoropyrimidin-4-ylamino)-2-methoxy-3-sulfonatophenylazo)-2-oxidophenylazo)-2,5,7-trisulfonato-4-naphtholato)copper(II); sodium/potassium (3-(4-(5-(5-chloro-4,6-difluoropyrimidin-2-ylamino)-2-methoxy-3-sulfonatophenylazo)-2-oxidophenylazo)-2,5,7-trisulfonato-4-naphtholato)copper(II)	407-100-7	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-075-00-X	reaction mass of: tris(3,5,5-trimethylhexylammonium) 4-amino-3-(4-(4-(2-amino-4-hydroxyphenylazo)anilino)-3-sulfonatophenylazo)-5,6-dihydro-5-oxo-6-phenylhydrazononaphthalene-2,7-disulfonate; tris(3,5,5-trimethylhexylammonium) 4-amino-3-(4-(4-(4-amino-2-hydroxyphenylazo)anilino)-3-sulfonatophenylazo)-5,6-dihydro-5-oxo-6-phenylhydrazononaphthalene-2,7-disulfonate (2:1)	406-000-0	—	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			
611-076-00-5	3-(2,6-dichloro-4-nitrophenylazo)-1-methyl-2-phenylindole	406-280-4	117584-16-4	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
611-077-00-0	dilithium disodium (5,5'-diamino-(μ-4,4'-dihydroxy-1:2-κ-2,O4,O4',-3,3'-[3,3'-dihydroxy-1:2-κ-2-O3,O3'-biphenyl-4,4'-ylenebisazo-1:2-(N3,N4-η:N3',N4'-η)]-dinaphthalene-2,7-disulfonato(8)))dicuprate(2-)	407-230-4	126637-70-5	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			
611-078-00-6	(2,2'-(3,3'-dioxidobiphenyl-4,4'-diyldiazo)bis(6-(4-(3-(diethylamino)propylamino)-6-(3-(diethylammonio)propylamino)-1,3,5-triazin-2-ylamino)-3-sulfonato-1-naphtholato))dicopper(II) acetate lactate	407-240-9	159604-94-1	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-079-00-1	disodium 7-[4-chloro-6-(N-ethyl-o-toluidino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(4-methoxy-2-sulfonatophenylazo)-2-naphthalenesulfonate	410-390-8	147703-64-8	Eye Dam. 1	H318	GHS05 Dgr	H318			
611-080-00-7	sodium 3-(2-acetamido-4-(4-(2-hydroxybutoxy)phenylazo)phenylazo)benzenesulfonate	410-150-2	147703-65-9	Skin Sens. 1	H317	GHS07 Wng	H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-081-00-2	tetrasodium [7-(2,5-dihydroxy-KO2-7-sulfonato-6-[4-(2,5,6-trichloro-pyrimidin-4-ylamino)phenylazo]-(N1,N7-N)-1-naphthylazo)-8-hydroxy-KO8-naphthalene-1,3,5-trisulfonato(6-)]cuprate(II)	411-470-5	141048-13-7	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
611-082-00-8	reaction mass of: pentasodium bis(1-(3(or 5)-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato)ferrate(1-); pentasodium [(1-(3-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato)-(5-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato]ferrate(1-)	407-570-3	—	Aquatic Chronic 2	H411	GHS09	H411			
611-083-00-3	reaction mass of: 2-[*N*-ethyl-4-[(5,6-dichlorobenzothiazol-2-yl)azo]-*m*-toludino]ethyl acetate; 2-[N-ethyl-4-[(6,7-dichlorobenzothiazol-2-yl)azo]-*m*-toludino]ethyl acetate (1:1)	411-560-4	—	STOT RE 1 Skin Sens. 1 Aquatic Chronic 2	H372 ** H317 H411	GHS08 GHS07 GHS09 Dgr	H372 ** H317 H411			
611-084-00-9	reaction mass of: N-(4-chlorophenyl)-4-(2,5-dichloro-4-(dimethylsulfamoyl)phenylazo)-3-hydroxy-2-naphthalenecarboxamide; N-(4-chlorophenyl)-4-(2,5-dichloro-4-(methylsulfamoyl)phenylazo)-3-hydroxy-2-naphthalenecarboxamide	412-550-2	—	Aquatic Chronic 4	H413	—	H413			
611-085-00-4	reaction mass of: 3-cyano-5-(2-cyano-4-nitro-phenylazo)-2-(2-hydroxy-ethylamino)-4-methyl-6-[3-(2-phenoxyethoxy)propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-6-(2-hydroxy-ethylamino)-4-methyl-2-[3-(2-phenoxyethoxy)propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-2-amino-4-methyl-6-[3-(3-hydroxypropoxy) propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-6-amino-4-methyl-2-[3-(3-methoxypropoxy) propylamino]pyridine	411-880-4	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-086-00-X	monolithium 5-[[2,4-dihydroxy-5-[(2-hydroxy-3,5-dinitrophenyl)azo]phenyl]azo]-2-naphthalenesulfonate], iron complex, monohydrate	411-360-7	—	Aquatic Chronic 3	H412	—	H412			
611-087-00-5	reaction mass of: 3-((5-cyano-1,6-dihydro-1,4-dimethyl-2-hydroxyl-6-oxo-3-pyridinyl) azo)-benzoyloxy-2-phenoxyethane; 3-((5-cyano-1,6-dihydro-1,4-dimethyl-2-hydroxy-6-oxo-3-pyridinyl)azo)-benzoyloxy-2-ethyloxy-2-(ethylphenol)	411-710-9	—	Aquatic Chronic 4	H413	—	H413			
611-088-00-0	reaction mass of: trilithium 4-amino-3-((4-((4-((2-amino-4-hydroxyphenyl)azo)phenyl) amino)-3-sulfophenyl)azo)-5-hydroxy-6-(phenylazo)naphthalene-2,7-disulfonate; trilithium 4-amino-3-((4-((4-((4-amino-2-hydroxyphenyl)azo)phenyl)amino)-3-sulfophenyl)azo)-5-hydroxy-6-(phenylazo) naphthalene-2,7-disulfonate	411-890-9	—	Acute Tox. 4 * Eye Dam. 1 Aquatic Chronic 3	H302 H318 H412	GHS05 GHS07 Dgr	H302 H318 H412			
611-089-00-6	2-((4-(ethyl-(2-hydroxyethyl)amino)-2-methylphenyl)azo)-6-methoxy-3-methyl-benzothiazolium methylsulfate	411-100-2	136213-73-5	STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H317 H400 H410	GHS08 GHS07 GHS09 Wng	H373 ** H317 H410			
611-090-00-1	2,5-dibutoxy-4-(morpholin-4-yl)benzene-diazonium 4-methylbenzenesulfonate	413-290-2	93672-52-7	Self-react. C Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H242 H302 H318 H317 H412	GHS02 GHS05 GHS07 Dgr	H242 H302 H318 H317 H412			T
611-091-00-7	sodium (1,0-1,95)/lithium (0,05-1) 5-((5-((5-chloro-6-fluoro-pyrimidin-4-yl)amino)-2-sulfonatophenyl)azo)-1,2-dihydro-6-hydroxy-1,4-dimethyl-2-oxo-3-pyridine-methylsulfonate	413-470-0	134595-59-8	Skin Sens. 1	H317	GHS07 Wng	H317			

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611-092-00-2	*tert*-(dodecyl/tetradecyl)-ammonium bis(3-(4-((5-(1,1-dimethyl-propyl)-2-hydroxy-3-nitrophenyl)azo)-3-methyl-5-hydroxy-(1*H*) pyrazol-1-yl)benzenesulfonamidato)chromate	413-210-6	—	Aquatic Chronic 2	H411	GHS09	H411			
611-093-00-8	sodium 2-(4-(4-fluoro-6-(2-sulfo-ethylamino)-[1,3,5]triazin-2-ylamino)-2-ureido-phenylazo)-5-(4-sulfophenylazo)benzene-1-sulfonate	410-770-3	146177-84-6	Skin Sens. 1	H317	GHS07 Wng	H317			
611-094-00-3	reaction mass of: 2-[2-acetylamino-4-[*N*,*N*-bis[2-ethoxy-carbonyloxy)ethyl]amino]phenylazo]-5,6-dichloro-1,3-benzothiazole; 2-[2-acetylamino-4-[*N*,*N*-bis[2-ethoxy-carbonyloxy)ethyl]amino]phenylazo]-6,7-dichloro-1,3-benzotriazole (1:1)	411-600-0	143145-93-1	Aquatic Chronic 4	H413	—	H413			
611-095-00-9	hexasodium 1,1'-[(1-amino-8-hydroxy-3,6-disulfonate-2,7-naphthalenediyl)bis(azo(4-sulfonate-1,3-phenyl)imino[6-[(4-chloro-3-sulfonatophenyl)amino]-1,3,5-triazin-2,4-diyl]]]bis[3-carboxypyridinium] dihydroxide	412-240-7	89797-03-5	Aquatic Chronic 2	H411	GHS09	H411			
611-096-00-4	methyl *N*-[3-acetylamino)-4-(2-cyano-4-nitrophenylazo)phenyl]-*N*-[(1-methoxy) acetyl]glycinate	413-040-2	149850-30-6	Skin Sens. 1	H317	GHS07 Wng	H317			
611-097-00-X	reaction mass of iron complexes of: 1,3-dihydroxy-4-[(5-phenylaminosulfonyl)-2-hydroxyphenylazo]-*n*-(5-amino-sulfonyl-2-hydroxyphenylazo)benzene and: 1,3-dihydroxy-4-[(5-phenylaminosulfonyl)-2-hydroxyphenylazo]-*n*-[4-(4-nitro-2-sulfophenylamino)phenylazo]benzene (*n*=2,5,6)	414-150-3	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-098-00-5	tetrakis(tetramethylammonium)3,3'-(6-(2-hydroxyethylamino)1,3,5-triazine-2,4-diyl-bisimino(2-methyl-4,1-phenyleneazo))bis-naphthalene-1,5-disulfonate	405-950-3	131013-83-7	Acute Tox. 3 * Aquatic Chronic 3	H301 H412	GHS06 Dgr	H301 H412			
611-099-00-0	(methylenebis(4,1-phenylenazo(1-(3-(dimethylamino)propyl)-1,2-dihydro-6-hydroxy-4-methyl-2-oxopyridine-5,3-diyl)))-1,1'-dipyridinium dichloride dihydrochloride	401-500-5	118658-99-4	Carc. 1B Aquatic Chronic 2	H350 H411	GHS08 GHS09 Dgr	H350 H411			

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611-100-00-4	potassium sodium 3,3'-(3(or4)-methyl-1,2-phenylenebis(imino(6-chloro)-1,3,5-triazine-4,2-diylimino(2-acetamido-5-methoxy)-4,1-phenylenazo)dinaphthalene-1,5-disulfonate	403-810-6	140876-13-7	Eye Dam. 1	H318	GHS05 Dgr	H318			
611-101-00-X	2'-(4-chloro-3-cyano-5-formyl-2-thienyl) azo-5'-diethylaminoacetanilide	405-200-5	104366-25-8	Skin Sens. 1	H317	GHS07 Wng	H317			
611-103-00-0	trisodium (1-(3-carboxylato-2-oxido-5-sulfonatophenylazo)-5-hydroxy-7-sulfonatonaphthalen-2-amido)nickel(II)	407-110-1	—	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			

611-104-00-6	reaction mass of: trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2(or 4or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2or 6)-(4-(4-nitro-2-sulfonatoanilino)phenylazo) phenolato)ferrate(1-);
	trisodium bis(2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)ferrate(1-);
	trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2 (or 4 or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2 or 6)-(4-nitro-2-sulfonatophenylazo)phenolato)ferrate(1-);
	trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2 (or 4 or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2 or 6)-(3-sulfonatophenylazo)phenolato)ferrate(1-);
	disodium 3,3'-(2,4-dihydroxy-1,3(or 1,5 or 3,5)-phenylenediazo)dibenzenesulfonate
		406-870-1	—	Skin Sens. 1
				Aquatic Chronic 2
					H317
					H411
						GHS07
						GHS09
						Wng
							H317
							H411			

611-105-00-1	sodium 4-(4-chloro-6-(*N*-ethylanilino)-1,3,5-triazin-2-ylamino)-2-(1-(2-chlorophenyl)-5-hydroxy-3-methyl-1*H*-pyrazol-4-ylazo)benzenesulfonate	407-800-2	136213-75-7	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			

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611-106-00-7	hexasodium 4,4'-dihydroxy-3,3'-bis[2-sulfonato-4-(4-sulfonatophenylazo)phenylazo]-7,7'[*p*-phenylenebis[imino(6-chloro-1,3,5-triazine-4,2-diyl)imino]]dinaphthalene-2-sulfonate	410-180-6	157627-99-1	Eye Dam. 1	H318	GHS05 Dgr	H318			
611-107-00-2	potassium sodium 4-(4-chloro-6-(3,6-disulfonato-7-(5,8-disulfonato-naphthalen-2-ylazo)-8-hydroxy-naphthalen-1-ylamino)-1,3,5-triazin-2-ylamino)-5-hydroxy-6-(4-(2-sulfatoethanesulfonyl)-phenylazo)-naphthalene-1,7-disulfonate	412-490-7	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-108-00-8	disodium 5-((4-((4-chloro-3-sulfonatophenyl)azo)-1-naphthyl)azo)-8-(phenylamino)-1-naphthalenesulfonate	413-600-6	6527-62-4	Aquatic Chronic 3	H412	—	H412			
611-109-00-3	reaction products of: copper(II) sulfate and tetrasodium 2,4-bis[6-(2-methoxy-5-sulfonatophenylazo)-5-hydroxy-7-sulfonato-2-naphthylamino]-6-(2-hydroxyethylamino)-1,3,5-triazine (2:1)	407-710-3	—	Aquatic Chronic 2	H411	GHS09	H411			
611-110-00-9	tetra-sodium/lithium 4,4'-bis-(8-amino-3,6-disulfonato-1-naphthol-2-ylazo)-3-methylazobenzene	408-210-8	124605-82-9	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-111-00-4	disodium 2-[[4-(2-chloroethylsulfonyl)phenyl]-[(2-hydroxy-5-sulfo-3-[3-[2-(2-(sulfooxy)ethylsulfonyl)ethylazo]-4-sulfobenzoato(3-)cuprate(1-)	414-230-8	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-112-00-X	tetrasodium 4-hydroxy-5-[4-[3-(2-sulfatoethanesulfonyl)phenylamino]-6-morpholin-4-yl-1,3,5-triazin-2-ylamino]-3-(1-sulfonatonaphthalen-2-ylazo)naphthalene-2,7-disulfonate	413-070-6	—	Skin Sens. 1	H317	GHS07 Wng	H317			
611-113-00-5	lithium sodium (2-(((5-((2,5-dichlorophenyl)azo)-2-hydroxyphenyl)methylene)amino) benzoato(2-))(2-((4,5-dihydro-3-methyl-5-oxo-1-phenyl-1*H*-pyrazol-4-yl)azo)-5-sulfobenzoato(3-)) chromate(2-)	414-280-0	149626-00-6	Aquatic Chronic 2	H411	GHS09	H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-114-00-0	lithium sodium (4-((5-chloro-2-hydroxyphenyl)azo)-2,4-dihydro-5-methyl-3*H*-pyrazol-3-onato(2-))(3-((4,5-dihydro-3-methyl-1-(4-methylphenyl)-5-oxo-1*H*-pyrazol-4-yl)azo)-4-hydroxy-5-nitrobenzenesulfonato(3-)) chromate(2-)	414-250-7	149564-66-9	Acute Tox. 4 * Eye Dam. 1 Aquatic Chronic 3	H302 H318 H412	GHS05 GHS07 Dgr	H302 H318 H412			
611-115-00-6	trilithium bis(4-((4-(diethylamino)-2-hydroxyphenyl)azo)-3-hydroxy-1-naphthalenesulfonato(3-))chromate(3-)	414-290-5	149564-65-8	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
611-116-00-1	reaction mass of: trisodium 5-{4-chloro-6-[2-(2,6-dichloro-5-cyanopyrimidin-4-ylamino)-propylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalene-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(2,6-dichloro-5-cyanopyrimidin-4-ylamino)-1-methyl-ethylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalene-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(4,6-dichloro-5-cyanopyrimidin-2-ylamino)-propylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(4,6-dichloro-5-cyanopyrimidin-2-ylamino)-1-methyl-ethylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate	414-620-8	—	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
611-117-00-7	1,3-bis{6-fluoro-4-[1,5-disulfo-4-(3-aminocarbonyl-1-ethyl-6-hydroxy-4-methyl-pyrid-2-on-5-ylazo)-phenyl-2-ylamino]-1,3,5-triazin-2-ylamino}propane lithium-, sodium salt	415-100-3	149850-29-3	Skin Sens. 1	H317	GHS07 Wng	H317			
611-118-00-2	sodium 1,2-bis[4-[4-{4-(4-sulfophenylazo)-2-sulfophenylazo}-2-ureido-phenyl-amino]-6-fluoro-1,3,5-triazin-2-ylamino]-propane, sodium salt	413-990-8		Skin Sens. 1	H317	GHS07 Wng	H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-119-00-8	tetrasodium 4-[4-chloro-6-(4-methyl-2-sulfophenylamino)-1,3,5-triazin-2-ylamino]-6-(4,5-dimethyl-2-sulfophenylazo)-5-hydroxy-naphthalene-2,7-disulfonate	415-400-4	148878-22-2	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
611-120-00-3	5-{4-[5-amino-2-[4-(2-sulfoxyethylsulfonyl) phenylazo]-4-sulfo-phenylamino]-6-chloro-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfo-naphthalen-2-ylazo)-naphthalene-2,7-disulfonicacid sodium salt	418-340-7	157707-94-3	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			
611-121-00-9	main component 6 (isomer): asym. 1:2 Cr(III)-complex of: A: 3-hydroxy-4-(2-hydroxy-naphthalene-1-ylazo) naphthalene-1-sulfonic acid, Na-salt and B: 1-[2-hydroxy-5-(4-methoxy-phenylazo)phenylazo]naphthalene-2-ol; main component 8 (isomer): asym. 1:2 Cr-complex of: A: 3-hydroxy-4-(2-hydroxy-naphthalene-1-ylazo)-naphthalene-1-sulfonic acid, Na-salt and B: 1-[2-hydroxy-5-(4-methoxy-phenylazo)-phenylazo]-naphthalene-2-ol	417-280-9	30785-74-1	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
611-122-00-4	hexasodium (di[*N*-(3-(4-[5-(5-amino-3-methyl-1-phenylpyrazol-4-yl-azo)-2,4-disulfo-anilino]-6-chloro-1,3,5-triazin-2-ylamino)phenyl)-sulfamoyl](di-sulfo)-phthalocyaninato)nickel	417-250-5	151436-99-6	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
611-123-00-X	3-(2,4-bis(4-((5-(4,6-bis(2-aminopropylamino)-1,3,5-triazin-2-ylamino)-4-hydroxy-2,7-disulfonaphthalen-3-yl)azo)phenylamino)-1,3,5-triazin-6-ylamino)propyldiethylammonium lactate	424-310-4	178452-66-9	Eye Dam. 1	H318	GHS05 Dgr	H318			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-124-00-5	reaction mass of: pentasodium 5-amino-3-(5-{4-chloro-6-[4-(2-sulfoxyethoxysulfonato)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo)-6-[5-(2,3-dibromopropionylamino)-2-sulfonatophenylazo]-4-hydroxynaphthalene-2,7-disulfonate; pentasodium 5-amino-6-[5-(2-bromoacryloylamino)-2-sulfonatophenylazo]-3-(5-{4-chloro-6-[4-(2-sulfoxyethoxysulfonato)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo)-4-hydroxynaphthalene-2,7-disulfonate; tetrasodium 5-amino-3-[5-{4-chloro-6-[4-(vinylsulfonyl)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo]-6-[5-(2,3-dibromopropionylamino)-2-sulfonatophenylazo]-4-hydroxynaphthalene-2,7-disulfonate	424-320-9		Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			
611-125-00-0	reaction mass of: Disodium 6-[3-carboxy-4,5-dihydro-5-oxo-4-sulfonatophenyl)pyrazolin-4-yl-azo]-3-[2-oxido-4-(ethensulfonyl)-5-methoxyphenylazo]-4-oxidonaphthalene-2-sulfonate copper (II) complex; Disodium 6-[3-carboxy-4,5-dihydro-5-oxo-4-sulfonatophenyl)pyrazolin-4-yl-azo]-3-[2-oxido-4-(2-hydroxyethylsulfonyl)-5-methoxyphenylazo]-4-oxidonaphthalene-2-sulfonate copper (II) complex	423-940-7	—	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			
611-126-00-6	2,6-bis-(2-(4-(4-amino-phenylamino)-phenylazo)-1,3-dimethyl-3*H*-imidazolium)-4-dimethylamino-1,3,5-triazine, dichloride	424-120-1	174514-06-8	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
611-127-00-1	pentasodium 4-amino-6-(5-(4-(2-ethyl-phenylamino)-6-(2-sulfatoethanesulfonyl)-1,3,5-triazin-2-ylamino)-2-sulfonatophenylazo)-5-hydroxy-3-(4-(2-sulfatoethanesulfonyl)phenylazo)naphthalene-2,7-disulfonate	423-790-2	—	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H318 H317 H412	GHS05 GHS07 Dgr	H318 H317 H412			G

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-128-00-7	*N,N'*-bis{6-chloro-4-[6-(4-vinylsulfonylphenylazo)-2,7-disulfonicacid-5-hydroxynapht-4-ylamino]-1,3,5-triazin-2-yl}-*N*-(2-hydroxyethyl)ethane-1,2-diamine, sodium salt	419-500-9	171599-85-2	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			
611-129-00-2	reaction mass of: 5-[(4-[(7-amino-1-hydroxy-3-sulfo-2-naphthyl)azo]-2,5-diethoxyphenyl)azo]-2-[(3-phosphonophenyl)azo]benzoic acid; 5-[(4-[(7-amino-1-hydroxy-3-sulfo-2-naphthyl)azo]-2,5-diethoxyphenyl)azo]-3-[(3-phosphonophenyl)azo]benzoic acid	418-230-9	163879-69-4	Expl. 1.3 **** Repr. 2 STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H203 H361f *** H373 ** H317 H411	GHS01 GHS08 GHS07 GHS09 Dgr	H203 H361f *** H373 ** H317 H411			
611-130-00-8	tetra-ammonium 2-[6-[7-(2-carboxylatophenylazo)-8-hydroxy-3,6-disulfonato-1-naphthylamino]-4-hydroxy-1,3,5-triazin-2-ylamino]benzoate	418-520-5	183130-96-3	Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H319 H400 H410	GHS07 GHS09 Wng	H319 H410			
611-131-00-3	2-[2-hydroxy-3-(2-chlorophenyl)carbamoyl-1-naphthylazo]-7-[2-hydroxy-3-(3-methylphenyl)carbamoyl-1-naphthylazo]fluoren-9-one	420-580-2	151798-26-4	Repr. 1B Aquatic Chronic 4	H360-D *** H413	GHS08 Dgr	H360D *** H413			
611-132-00-9	pentasodium bis{7-[4-(1-butyl-5-cyano-1,2-dihydro-2-hydroxy-4-methyl-6-oxo-3-pyridylazo)phenylsulfonylamino]-5'-nitro-3,3'-disulfonatonaphthalene-2-azobenzene-1,2'-diolato} chromate (III)	419-210-2	178452-71-6	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			
611-133-00-4	Product by process iron complex of azo dyestuffs obtained by coupling a mixture of diazotized 2-amino-1-hydroxybenzene-4-sulfanilide and 2-amino-1-hydroxybenzene-4-sulfonamide with resorcin, the obtained mixture being subsequently submitted to a second coupling reaction with a mixture of diazotized 3-aminobenzene-1-sulfonic acid (metanilic acid) and 4'-amino-4-nitro-1,1'-diphenylamine-2-sulfonic acid and metallization with ferric chloride, sodium salt	419-260-5	—	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
611-134-00-X	trisodium 2-{α[2-hydroxy-3-[4-chloro-6-[4-(2,3-dibromopropionylamino)-2-sulfonatophenylamino]-1,3,5-triazin-2-ylamino]-5-sulfonatophenylazo]-benzylidenehydrazino}-4-sulfonatobenzoate, copper complex	423-770-3	—	Eye Dam. 1 Aquatic Chronic 2	H318 H411	GHS05 GHS09 Dgr	H318 H411			
611-135-00-5	reaction product of: 2-[[4-amino-2-ureidophenylazo]-5-[(2-(sulfooxy)ethyl)sulfonyl]] benzenesulfonic acid with 2,4,6-trifluoropyrimidine and partial hydrolysis to the corresponding vinylsulfonyl derivative, mixed potassium/sodium salt	424-250-9	—	Eye Dam. 1 Aquatic Chronic 3	H318 H412	GHS05 Dgr	H318 H412			
611-136-00-0	2-{4-(2-ammoniopropylamino)-6-[4-hydroxy-3-(5-methyl-2-methoxy-4-sulfamoylphenylazo)-2-sulfonatonaphth-7-ylamino]-1,3,5-triazin-2-ylamino}-2-aminopropyl formate	424-260-3	—	Repr. 2 Eye Dam. 1 Aquatic Chronic 2	H361f *** H318 H411	GHS05 GHS08 GHS09 Dgr	H361f *** H318 H411			
611-137-00-6	6-*tert*-butyl-7-chloro-3-tridecyl-7,7a-dihydro-1*H*-pyrazolo[5,1-c]-1,2,4-triazole	419-870-1	159038-16-1	Aquatic Chronic 4	H413	—	H413			
611-138-00-1	2-(4-aminophenyl)-6-*tert*-butyl-1*H*-pyrazolo [1,5-b][1,2,4]triazole	415-910-7	152828-25-6	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
611-140-00-2	azafenidin (ISO); 2-(2,4-dichloro-5-prop-2-ynyloxyphenyl)-5,6,7,8-tetrahydro-1,2,4-triazolo[4,3-a]pyridin-3(2*H*)-one	—	68049-83-2	Repr. 1B STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H360-Df H373 ** H400 H410	GHS08 GHS09 Dgr	H360Df H373 ** H410		M = 1 000	
612-001-00-9	mono-methylamine; [1] di-methylamine; [2] tri-methylamine [3]	200-820-0 [1] 204-697-4 [2] 200-875-0 [3]	74-89-5 [1] 124-40-3 [2] 75-50-3 [3]	Flam. Gas 1 Press. Gas Acute Tox. 4 * STOT SE 3 Skin Irrit. 2 Eye Dam. 1	H220 H332 H335 H315 H318	GHS02 GHS04 GHS05 GHS07 Dgr	H220 H332 H335 H315 H318		* Skin Irrit. 2; H315: C ≥ 5 % Eye Dam. 1; H318: C ≥ 5 % Eye Irrit. 2; H319: 0,5 % ≤ C < 5 % STOT SE 3; H335: C ≥ 5 %	U 5

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612-001-01-6	mono-methylamine ... %; [1] di-methylamine ... %; [2] tri-methylamine ... % [3]	200-820-0 [1] 204-697-4 [2] 200-875-0 [3]	74-89-5 [1] 124-40-3 [2] 75-50-3 [3]	Flam. Liq. 1 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H224 H332 H302 H314	GHS02 GHS05 GHS07 Dgr	H224 H332 H302 H314		* STOT SE 3; H335: C ≥ 5 %	B
612-002-00-4	ethylamine	200-834-7	75-04-7	Flam. Gas 1 Press. Gas Eye Irrit. 2 STOT SE 3	H220 H319 H335	GHS02 GHS04 GHS07 Dgr	H220 H319 H335			U
612-003-00-X	diethylamine	203-716-3	109-89-7	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H225 H332 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H312 H302 H314		STOT SE 3; H335: C ≥ 1 %	
612-004-00-5	triethylamine	204-469-4	121-44-8	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H225 H332 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H312 H302 H314		STOT SE 3; H335: C ≥ 1 %	
612-005-00-0	butylamine	203-699-2	109-73-9	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H225 H332 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H312 H302 H314		STOT SE 3; H335: C ≥ 1 %	
612-006-00-6	ethylenediamine; 1,2-diaminoethane	203-468-6	107-15-3	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Resp. Sens. 1 Skin Sens. 1	H226 H312 H302 H314 H334 H317	GHS02 GHS08 GHS05 GHS07 Dgr	H226 H312 H302 H314 H334 H317			

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612-007-00-1	2-aminopropane; isopropylamine	200-860-9	75-31-0	Flam. Liq. 1 Eye Irrit. 2 STOT SE 3 Skin Irrit. 2	H224 H319 H335 H315	GHS02 GHS07 Dgr	H224 H319 H335 H315			
612-008-00-7	aniline	200-539-3	62-53-3	Carc. 2 Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H351 H341 H331 H311 H301 H372 ** H318 H317 H400	GHS06 GHS08 GHS05 GHS09 Dgr	H351 H341 H331 H311 H301 H372 ** H318 H317 H400		* STOT RE 1; H372: C ≥ 1 % STOT RE 2; H373: 0,2 % ≤ C < 1 %	
612-009-00-2	salts of aniline	—	—	Carc. 2 Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H351 H341 H331 H311 H301 H372 ** H318 H317 H400	GHS06 GHS08 GHS05 GHS09 Dgr	H351 H341 H331 H311 H301 H372 ** H318 H317 H400		* STOT RE 1; H372: C ≥ 1 % STOT RE 2; H373: 0,2 % ≤ C < 1 %	A
612-010-00-8	chloroanilines (with exception of those specified elsewhere in this Annex)	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			C

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-011-00-3	4-nitrosoaniline	211-535-6	659-49-4	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H332 H312 H302	GHS07 Wng	H332 H312 H302			
612-012-00-9	*o*-nitroaniline; [1] *m*-nitroaniline; [2] *p*-nitroaniline [3]	201-855-4 [1] 202-729-1 [2] 202-810-1 [3]	88-74-4 [1] 99-09-2 [2] 100-01-6 [3]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 3	H331 H311 H301 H373 ** H412	GHS06 GHS08 Dgr	H331 H311 H301 H373 ** H412			C
612-013-00-4	3-aminobenzene sulphonic acid; metanilic acid	204-473-6	121-47-1	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H332 H312 H302	GHS07 Wng	H332 H312 H302			
612-014-00-X	sulphanilic acid; 4-aminobenzenesulphonic acid	204-482-5	121-57-3	Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1	H319 H315 H317	GHS07 Wng	H319 H315 H317			
612-015-00-5	*N*-methylaniline	202-870-9	100-61-8	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			
612-016-00-0	*N*,*N*-dimethylaniline	204-493-5	121-69-7	Carc. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H351 H331 H311 H301 H411	GHS06 GHS08 GHS09 Dgr	H351 H331 H311 H301 H411			

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612-017-00-6	*N*-methyl-*N*,2,4,6-tetranitroaniline; tetryl	207-531-9	479-45-8	Expl. 1.1 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H201 H331 H311 H301 H373 **	GHS01 GHS06 GHS08 Dgr	H201 H331 H311 H301 H373 **			
612-018-00-1	bis(2,4,6-trinitrophenyl)amine; hexyl	205-037-8	131-73-7	Expl. 1.1 Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Chronic 2	H201 H330 H310 H300 H373 ** H411	GHS01 GHS06 GHS08 GHS09 Dgr	H201 H330 H310 H300 H373 ** H411			
612-019-00-7	dipicrylamine, ammonium salt	220-639-0	2844-92-0	Expl. 1.1 Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Chronic 2	H201 H330 H310 H300 H373 ** H411	GHS01 GHS06 GHS08 GHS09 Dgr	H201 H330 H310 H300 H373 ** H411	EUH001		
612-020-00-2	1-naphthylamine	205-138-7	134-32-7	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
612-022-00-3	2-naphthylamine	202-080-4	91-59-8	Carc. 1A Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411		Carc. 1A; H350: C ≥ 0,01 %	

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-023-00-9	phenylhydrazine; [1] phenylhydrazinium chloride; [2] phenylhydrazine hydrochloride; [3] phenylhydrazinium sulphate (2:1) [4]	202-873-5 [1] 200-444-7 [2] 248-259-0 [3] 257-622-2 [4]	100-63-0 [1] 59-88-1 [2] 27140-08-5 [3] 52033-74-6 [4]	Carc. 1B Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1	H350 H341 H331 H311 H301 H372 ** H319 H315 H317 H400	GHS06 GHS08 GHS09 Dgr	H350 H341 H331 H311 H301 H372 ** H319 H315 H317 H400			
612-024-00-4	*m*-toluidine; 3-aminotoluene	203-583-1	108-44-1	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1	H331 H311 H301 H373 ** H400	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H400			
612-025-00-X	nitrotoluidines, with the exception of those specified elsewhere in this Annex	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411			C
612-026-00-5	diphenylamine	204-539-4	122-39-4	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H373 ** H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-027-00-0	xylidines with the exception of those specified elsewhere in this Annex; dimethyl anilines with the exception of those specified elsewhere in this Annex	—	—	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411			C
612-028-00-6	*p*-phenylenediamine	203-404-7	106-50-3	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H319 H317 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H319 H317 H410			
612-029-00-1	benzene-1,4-diamine dihydrochloride; *p*-phenylenediamine dihydrochloride	210-834-9	624-18-0	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H319 H317 H400 H410	GHS06 GHS09 Dgr	H331 H311 H301 H319 H317 H410			
612-030-00-7	2-methyl-*p*-phenylenediamine sulphate [1]	210-431-8 [1] 228-871-4 [2]	615-50-9 [1] 6369-59-1 [2]	Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H301 H332 H312 H317 H411	GHS06 GHS09 Dgr	H301 H332 H312 H317 H411			
612-031-00-2	*N*,*N*-dimethylbenzene-1,3-diamine; [1] 4-amino-*N*,*N*-dimethylaniline; 3-amino-*N*,*N*'-dimethylaniline [2]	220-623-3 [1] 202-807-5 [2]	2836-04-6 [1] 99-98-9 [2]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H331 H311 H301	GHS06 Dgr	H331 H311 H301			C

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-032-00-8	*N*,*N*,*N*',*N*'-tetramethyl-*p*-phenylenediamine	202-831-6	100-22-1	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H332 H312 H302	GHS07 Wng	H332 H312 H302			
612-033-00-3	2-aminophenol	202-431-1	95-55-6	Muta. 2 Acute Tox. 4 * Acute Tox. 4 *	H341 H332 H302	GHS08 GHS07 Wng	H341 H332 H302			
612-034-00-9	2-amino-4,6-dinitrophenol; picramic acid (< 20 % water)	202-544-6	96-91-3	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H201 H332 H312 H302 H412	GHS01 GHS07 Wng	H201 H332 H312 H302 H412			T
612-034-01-6	2-amino-4,6-dinitrophenol; picramic acid; [≥ 20 % water]	202-544-6	96-91-3	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H332 H312 H302 H412	GHS07 Wng	H332 H312 H302 H412			G
612-035-00-4	2-methoxyaniline; *o*-anisidine	201-963-1	90-04-0	Carc. 1B Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 *	H350 H341 H331 H311 H301	GHS06 GHS08 Dgr	H350 H341 H331 H311 H301			
612-036-00-X	3,3'-dimethoxybenzidine; *o*-dianisidine	204-355-4	119-90-4	Carc. 1B Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			
612-037-00-5	salts of 3,3'-dimethoxybenzidine; salts of *o*-dianisidine	—	—	Carc. 1B Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			A

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-038-00-0	2-nitro-*p*-anisidine; 4-methoxy-2-nitroaniline	202-547-2	96-96-8	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Chronic 3	H330 H310 H300 H373 ** H412	GHS06 GHS08 Dgr	H330 H310 H300 H373 ** H412			
612-039-00-6	2-ethoxyaniline; *o*-phenetidine	202-356-4	94-70-2	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H331 H311 H301 H373 **	GHS06 GHS08 Dgr	H331 H311 H301 H373 **			
612-040-00-1	2,4-dinitroaniline	202-553-5	97-02-9	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Chronic 2	H330 H310 H300 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H330 H310 H300 H373 ** H411			
612-041-00-7	4,4'-bi-*o*-toluidine	204-358-0	119-93-7	Carc. 1B Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411			
612-042-00-2	benzidine; 1,1'-biphenyl-4,4'-diamine; 4,4'-diaminobiphenyl; biphenyl-4,4'-ylenediamine	202-199-1	92-87-5	Carc. 1A Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H302 H410		Carc. 1A; H350: C ≥ 0,01 %	
612-043-00-8	*N,N'*-dimethylbenzidine	—	2810-74-4	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H332 H312 H302	GHS07 Wng	H332 H312 H302			
612-044-00-3	*N,N'*-diacetylbenzidine	210-338-2	613-35-4	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H332 H312 H302	GHS07 Wng	H332 H312 H302			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-046-00-4	allylamine	203-463-9	107-11-9	Flam. Liq. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H225 H331 H311 H301 H411	GHS02 GHS06 GHS09 Dgr	H225 H331 H311 H301 H411			
612-047-00-X	benzylamine	202-854-1	100-46-9	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H312 H302 H314	GHS05 GHS07 Dgr	H312 H302 H314			
612-048-00-5	dipropylamine	205-565-9	142-84-7	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H225 H332 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H312 H302 H314		STOT SE 3; H335: C ≥ 1 %	
612-049-00-0	di-*n*-butylamine; [1] di-*sec*-butylamine [2]	203-921-8 [1] 210-937-9 [2]	111-92-2 [1] 626-23-3 [2]	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H226 H332 H312 H302	GHS02 GHS07 Wng	H226 H332 H312 H302			
612-050-00-6	cyclohexylamine	203-629-0	108-91-8	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H226 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H226 H312 H302 H314			
612-051-00-1	4,4'-diaminodiphenylmethane; 4,4'-methylenedianiline	202-974-4	101-77-9	Carc. 1B Muta. 2 STOT SE 1 STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H350 H341 H370 ** H373 ** H317 H411	GHS08 GHS07 GHS09 Dgr	H350 H341 H370 ** H373 ** H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-052-00-7	(*S*)-*sec*-butylamine; (*S*)-2-aminobutane; [1] (*R*)-*sec*-butylamine; (*R*)-2-aminobutane; [2] *sec*-butylamine; 2-aminobutane [3]	208-164-7 [1] 236-232-6 [2] 237-732-7 [3]	513-49-5 [1] 13250-12-9 [2] 13952-84-6 [3]	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A Aquatic Acute 1	H225 H332 H302 H314 H400	GHS02 GHS05 GHS07 GHS09 Dgr	H225 H332 H302 H314 H400			C
612-053-00-2	*N*-ethylaniline	203-135-5	103-69-5	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H331 H311 H301 H373 **	GHS06 GHS08 Dgr	H331 H311 H301 H373 **			
612-054-00-8	*N*,*N*-diethylaniline	202-088-8	91-66-7	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411		*	
612-055-00-3	*N*-methyl-*o*-toluidine; [1] *N*-methyl-*m*-toluidine; [2] *N*-methyl-*p*-toluidine [3]	210-260-9 [1] 211-795-0 [2] 210-769-6 [3]	611-21-2 [1] 696-44-6 [2] 623-08-5 [3]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 3	H331 H311 H301 H373 ** H412	GHS06 GHS08 Dgr	H331 H311 H301 H373 ** H412			C
612-056-00-9	*N*,*N*-dimethyl-*p*-toluidine; [1] *N*,*N*-dimethyl-*m*-toluidine; [2] *N*,*N*-dimethyl-*o*-toluidine [3]	202-805-4 [1] 204-495-6 [2] 210-199-8 [3]	99-97-8 [1] 121-72-2 [2] 609-72-3 [3]	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 3	H331 H311 H301 H373 ** H412	GHS06 GHS08 Dgr	H331 H311 H301 H373 ** H412		*	C

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-057-00-4	piperazine	203-808-3	110-85-0	Skin Corr. 1B Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 3	H314 H334 H317 H412	GHS08 GHS05 Dgr	H314 H334 H317 H412			
612-058-00-X	2,2'-iminodiethylamine; diethylenetriamine	203-865-4	111-40-0	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1	H312 H302 H314 H317	GHS05 GHS07 Dgr	H312 H302 H314 H317			
612-059-00-5	3,6-diazaoctanethylenediamin; triethylenetetramine	203-950-6	112-24-3	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 3	H312 H314 H317 H412	GHS05 GHS07 Dgr	H312 H314 H317 H412			
612-060-00-0	3,6,9-triazaundecamethylenediamine; tetraethylenepentamine	203-986-2	112-57-2	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 2	H312 H302 H314 H317 H411	GHS05 GHS07 GHS09 Dgr	H312 H302 H314 H317 H411			
612-061-00-6	3-aminopropyldimethylamine; *N*,*N*-dimethyl-1,3-diaminopropane	203-680-9	109-55-7	Flam. Liq. 3 Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1	H226 H302 H314 H317	GHS02 GHS05 GHS07 Dgr	H226 H302 H314 H317			
612-062-00-1	3-aminopropyldiethylamine; *N*,*N*-diethyl-1,3-diaminopropane	203-236-4	104-78-9	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1	H226 H312 H302 H314 H317	GHS02 GHS05 GHS07 Dgr	H226 H312 H302 H314 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-063-00-7	3,3'-iminodi(propylamine); dipropylenetriamine	200-261-2	56-18-8	Acute Tox. 2 * Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1A Skin Sens. 1	H330 H311 H302 H314 H317	GHS06 GHS05 Dgr	H330 H311 H302 H314 H317			
612-064-00-2	3,6,9,12-tetra-azatetradecamethylenediamine; pentacthylenehexamine	223-775-9	4067-16-7	Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H314 H317 H410			
612-065-00-8	polyethlyenepolyamines with the exception of those specified elsewhere in this Annex	—	—	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H312 H302 H314 H317 H410			
612-066-00-3	dicyclohexylamine	202-980-7	101-83-7	Acute Tox. 4 * Skin Corr. 1B Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H314 H410			
612-067-00-9	3-aminomethyl-3,5,5-trimethylcyclohexylamine	220-666-8	2855-13-2	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 3	H312 H302 H314 H317 H412	GHS05 GHS07 Dgr	H312 H302 H314 H317 H412			

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612-068-00-4	3,3'-dichlorobenzidine; 3,3'-dichlorobiphenyl-4,4'-ylenediamine	202-109-0	91-94-1	Carc. 1B Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H312 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H312 H317 H410			
612-069-00-X	salts of 3,3'-dichlorobenzidine; salts of 3,3'-dichlorobiphenyl-4,4'-ylenediamine	—	—	Carc. 1B Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H312 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H312 H317 H410			A
612-070-00-5	salts of benzidine	208-519-6 208-520-1 244-236-4 252-984-8	531-85-1 531-86-2 21136-70-9 36341-27-2	Carc. 1A Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H302 H410			A
612-071-00-0	salts of 2-naphthylamine	209-030-0 210-313-6	553-00-4 612-52-2	Carc. 1A Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411			A
612-072-00-6	biphenyl-4-ylamine; xenylamine; 4-aminobiphenyl	202-177-1	92-67-1	Carc. 1A Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			
612-073-00-1	salts of biphenyl-4-ylamine; salts of xenylamine; salts of 4-aminobiphenyl	—	—	Carc. 1A Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			A

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-074-00-7	benzyldimethylamine	203-149-1	103-83-3	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Aquatic Chronic 3	H226 H332 H312 H302 H314 H412	GHS02 GHS05 GHS07 Dgr	H226 H332 H312 H302 H314 H412			
612-075-00-2	2-aminoethyldimethylamine; 2-dimethylaminoethylamine	203-541-2	108-00-9	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H225 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H312 H302 H314			
612-076-00-8	ethyldimethylamine	209-940-8	598-56-1	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H225 H332 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H302 H314			
612-077-00-3	dimethylnitrosoamine; N-nitrosodimethylamine	200-549-8	62-75-9	Carc. 1B Acute Tox. 2 * Acute Tox. 3 * STOT RE 1 Aquatic Chronic 2	H350 H330 H301 H372 ** H411	GHS06 GHS08 GHS09 Dgr	H350 H330 H301 H372 ** H411		Carc. 1B; H350: C ≥ 0,001 %	
612-078-00-9	2,2'-dichloro-4,4'-methylenedianiline; 4,4'-methylene bis(2-chloroaniline)	202-918-9	101-14-4	Carc. 1B Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H302 H410			
612-079-00-4	salts of 2,2'-dichloro-4,4'-methylenedianiline; salts of 4,4'-methylenebis(2-chloroaniline)	—	—	Carc. 1B Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H302 H410			A

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-080-00-X	4-amino-*N*,*N*-diethylaniline; *N*,*N*-diethyl-p-phenylendiamine	202-214-1	93-05-0	Acute Tox. 3 * Skin Corr. 1B	H301 H314	GHS06 GHS05 Dgr	H301 H314			
612-081-00-5	salts of 4,4'-bi-*o*-toluidine; salts of 3,3'-dimethylbenzidine; salts of *o*-tolidine	210-322-5 265-294-7 277-985-0	612-82-8 64969-36-4 74753-18-7	Carc. 1B Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411			A
612-082-00-0	thiourea; thiocarbamide	200-543-5	62-56-6	Carc. 2 Repr. 2 Acute Tox. 4 * Aquatic Chronic 2	H351 H361d *** H302 H411	GHS08 GHS07 GHS09 Wng	H351 H361d *** H302 H411			
612-083-00-6	1-methyl-3-nitro-1-nitrosoguanidine	200-730-1	70-25-7	Carc. 1B Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 2	H350 H332 H319 H315 H411	GHS08 GHS07 GHS09 Dgr	H350 H332 H319 H315 H411		Carc. 1B; H350: C ≥ 0,01 %	
612-084-00-1	dapsone; 4,4'-diamino diphenyl sulfone	201-248-4	80-08-0	Acute Tox. 4 *	H302	GHS07 Wng	H302			
612-085-00-7	4,4'-methylenedi-*o*-toluidine	212-658-8	838-88-0	Carc. 1B Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H350 H302 H317 H410			
612-086-00-2	amitraz (ISO); *N*,*N*-bis(2,4-xylyliminomethyl) methylamine	251-375-4	33089-61-1	Acute Tox. 4 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H373 ** H317 H400 H410	GHS08 GHS07 GHS09 Wng	H302 H373 ** H317 H410		M = 10	

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-087-00-8	guazatine (ISO)		108173-90-6	Acute Tox. 2 * Acute Tox. 4 * Acute Tox. 4 * STOT SE 3 Skin Irrit. 2 Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H330 H312 H302 H335 H315 H318 H400 H410	GHS06 GHS05 GHS09 Dgr	H330 H312 H302 H335 H315 H318 H410			
612-088-00-3	simazine (ISO); 6-chloro-*N*,*N*'-diethyl-1,3,5-triazine-2,4-diamine	204-535-2	122-34-9	Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H400 H410	GHS08 GHS09 Wng	H351 H410			
612-089-00-9	1,5-naphthylenediamine	218-817-8	2243-62-1	Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H400 H410	GHS08 GHS09 Wng	H351 H410			
612-090-00-4	2,2'-(nitrosoimino)bisethanol	214-237-4	1116-54-7	Carc. 1B	H350	GHS08 Dgr	H350			
612-091-00-X	*o*-toluidine; 2-aminotoluene	202-429-0	95-53-4	Carc. 1B Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Aquatic Acute 1	H350 H331 H301 H319 H400	GHS06 GHS08 GHS09 Dgr	H350 H331 H301 H319 H400			
612-092-00-5	*N*,*N*'-(2,2-dimethylpropylidene)hexamethylenediamine	401-660-6	1000-78-8	Skin Irrit. 2 Skin Sens. 1	H315 H317	GHS07 Wng	H315 H317			
612-093-00-0	3,5-dichloro-4-(1,1,2,2-tetrafluoroethoxy) aniline	401-790-3	104147-32-2	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-094-00-6	4-(2-chloro-4-trifluoromethyl)phenoxy-2-fluoroaniline hydrochloride	402-190-4	—	STOT RE 1 Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H372 ** H302 H318 H317 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H372 ** H302 H318 H317 H410			
612-095-00-1	benzyl-2-hydroxydodecyldimethylammonium benzoate	402-610-6	113694-52-3	Skin Corr. 1B Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H314 H302 H400 H410	GHS05 GHS07 GHS09 Dgr	H314 H302 H410			
612-096-00-7	4,4'-carbonimidoylbis[*N*,*N*-dimethylaniline]	207-762-5	492-80-8	Carc. 2 Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 2	H351 H302 H319 H411	GHS08 GHS07 GHS09 Wng	H351 H302 H319 H411			
612-097-00-2	salts of 4,4'-carbonimidoylbis[*N*,*N*-dimethylaniline]	—	—	Carc. 2 Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 2	H351 H302 H319 H411	GHS08 GHS07 GHS09 Wng	H351 H302 H319 H411			A
612-098-00-8	nitrosodipropylamine	210-698-0	621-64-7	Carc. 1B Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411		Carc. 1B; H350: C ≥ 0,001 %	
612-099-00-3	4-methyl-*m*-phenylenediamine; 2,4-toluenediamine	202-453-1	95-80-7	Carc. 1B Acute Tox. 3 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H350 H301 H312 H319 H317 H411	GHS06 GHS08 GHS09 Dgr	H350 H301 H312 H319 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-100-00-7	propylenediamine	201-155-9	78-90-0	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A	H226 H312 H302 H314	GHS02 GHS05 GHS07 Dgr	H226 H312 H302 H314			
612-101-00-2	methenamine; hexamethylenetetramine	202-905-8	100-97-0	Flam. Sol. 2 Resp. Sens. 1 Skin Sens. 1	H228 H334 H317	GHS02 GHS08 Dgr	H228 H334 H317			
612-102-00-8	*N*,*N*-bis(3-aminopropyl)methylamine	203-336-8	105-83-9	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1B	H331 H311 H302 H314	GHS06 GHS05 Dgr	H331 H311 H302 H314			
612-103-00-3	*N*,*N*,*N*',*N*'-tetramethylethylenediamine	203-744-6	110-18-9	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H225 H332 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H302 H314			
612-104-00-9	hexamethylenediamine	204-679-6	124-09-4	Acute Tox. 4 * Acute Tox. 4 * STOT SE 3 Skin Corr. 1B	H312 H302 H335 H314	GHS05 GHS07 Dgr	H312 H302 H335 H314			
612-105-00-4	2-piperazin-1-ylethylamine	205-411-0	140-31-8	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 3	H312 H302 H314 H317 H412	GHS05 GHS07 Dgr	H312 H302 H314 H317 H412			
612-106-00-X	2,6-diethylaniline	209-445-7	579-66-8	Acute Tox. 4 *	H302	—	H302			
612-107-00-5	1-phenylethylamine; [1] DL-α-methylbenzylamine [2]	202-706-6 [1] 210-545-8 [2]	98-84-0 [1] 618-36-0 [2]	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H312 H302 H314	GHS05 GHS07 Dgr	H312 H302 H314			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-108-00-0	3-aminopropyltriethoxysilane	213-048-4	919-30-2	Acute Tox. 4 * Skin Corr. 1B	H302 H314	GHS05 GHS07 Dgr	H302 H314			
612-109-00-6	bis(2-dimethylaminoethyl)(methyl)amine	221-201-1	3030-47-5	Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1B	H311 H302 H314	GHS06 GHS05 Dgr	H311 H302 H314			
612-110-00-1	2,2'-dimethyl-4,4'-methylenebis(cyclohexylamine)	229-962-1	6864-37-5	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1A Aquatic Chronic 2	H331 H311 H302 H314 H411	GHS06 GHS05 GHS09 Dgr	H331 H311 H302 H314 H411			
612-111-00-7	2-methyl-*m*-phenylenediamine; 2,6-toluenediamine	212-513-9	823-40-5	Muta. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H341 H312 H302 H317 H411	GHS08 GHS07 GHS09 Wng	H341 H312 H302 H317 H411			
612-112-00-2	*p*-anisidine; 4-methoxyaniline	203-254-2	104-94-9	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * STOT RE 2 * Aquatic Acute 1	H330 H310 H300 H373 ** H400	GHS06 GHS08 GHS09 Dgr	H330 H310 H300 H373 ** H400			
612-113-00-8	6-methyl-2,4-bis(methylthio)phenylene-1,3-diamine	403-240-8	106264-79-3	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-114-00-3	*R*,*R*-2-hydroxy-5-(1-hydroxy-2-(4-phenylbut-2-ylamino)ethyl)benzamide hydrogen 2,3-bis(benzoyloxy)succinate	404-390-7	—	Flam. Sol. 1 Skin Sens. 1 Aquatic Chronic 3	H228 H317 H412	GHS02 GHS07 Wng	H228 H317 H412			
612-115-00-9	dimethyldioctadecylammonium hydrogen sulfate	404-050-8	123312-54-9	Eye Irrit. 2 Aquatic Chronic 4	H319 H413	GHS07 Wng	H319 H413			
612-116-00-4	C_{8-18}alkylbis(2-hydroxyethyl)ammonium bis(2-ethylhexyl)phosphate	404-690-8	68132-19-4	Acute Tox. 3 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H314 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H331 H314 H317 H410			
612-117-00-X	C_{12-14}-*tert*-alkylamine, methylphosphonic acid salt	404-750-3	119415-07-5	Acute Tox. 4 * Skin Corr. 1B Aquatic Chronic 2	H302 H314 H411	GHS05 GHS07 GHS09 Dgr	H302 H314 H411			
612-118-00-5	A reaction mass of: (1,3-dioxo-2*H*-benz(de)isoquinolin-2-ylpropyl)hexadecyldimethylammonium 4-toluenesulfonate; (1,3-dioxo-2*H*-benz(de)isoquinolin-2-ylpropyl)hexadecyldimethylammonium bromide	405-080-4	—	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
612-119-00-0	benzyldimethyloctadecylammonium 3-nitrobenzenesulfonate	405-330-2	—	Skin Irrit. 2 Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H315 H318 H400 H410	GHS05 GHS09 Dgr	H315 H318 H410			
612-120-00-6	aclonifen (ISO); 2-chloro-6-nitro-3-phenoxyaniline	277-704-1	74070-46-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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612-121-00-1	amines, polyethylenepoly-; HEPA	268-626-9	68131-73-7	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H312 H302 H314 H317 H410			
612-122-00-7	hydroxylamine	232-259-2	7803-49-8	Unst. Expl. Met. Corr. 1 Acute Tox. 4 * STOT RE 2 * STOT SE 3 Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H200 H290 H302 H373 ** H335 H315 H318 H317 H400	GHS01 GHS08 GHS05 GHS07 GHS09 Dgr	H200 H290 H302 H373 ** H335 H315 H318 H317 H400			T
612-123-00-2	hydroxylammonium chloride; hydroxylamine hydrochloride; [1] bis(hydroxylammonium) sulphate; hydroxylamine sulphate (2:1); [2] hydroxylammonium hydrogensulphate; hydroxylamine sulphate (1:1) [3]	226-798-2 [1] 233-118-8 [2] 233-154-4 [3]	5470-11-1 [1] 10039-54-0 [2] 10046-00-1 [3]	Met. Corr. 1 Acute Tox. 4 * STOT RE 2 * Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1	H290 H302 H373 ** H319 H315 H317 H400	GHS08 GHS05 GHS07 GHS09 Wng	H290 H302 H373 ** H319 H315 H317 H400			G T
612-124-00-8	*N*,*N*,*N*-trimethylanilinium chloride	205-319-0	138-24-9	Acute Tox. 3 * Acute Tox. 3 *	H311 H301	GHS06 Dgr	H311 H301			
612-125-00-3	2-methyl-*p*-phenylenediamine; 2,5-toluenediamine	202-442-1	95-70-5	Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H301 H332 H312 H317 H411	GHS06 GHS09 Dgr	H301 H332 H312 H317 H411			

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612-126-00-9	toluene-2,4-diammonium sulphate; 4-methyl-m-phenylenediamine sulfate	265-697-8	65321-67-7	Carc. 1B Acute Tox. 3 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H350 H301 H312 H319 H317 H411	GHS06 GHS08 GHS09 Dgr	H350 H301 H312 H319 H317 H411			
612-127-00-4	3-aminophenol	209-711-2	591-27-5	Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 2	H332 H302 H411	GHS07 GHS09 Wng	H332 H302 H411			
612-128-00-X	4-aminophenol	204-616-2	123-30-8	Muta. 2 Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H341 H332 H302 H400 H410	GHS08 GHS07 GHS09 Wng	H341 H332 H302 H410			
612-129-00-5	diisopropylamine	203-558-5	108-18-9	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H225 H332 H302 H314	GHS02 GHS05 GHS07 Dgr	H225 H332 H302 H314		STOT SE 3; H335: C ≥ 5 %	
612-130-00-0	2,6-diamino-3,5-diethyltoluene; 4,6-diethyl-2-methyl-1,3-benzenediamine; [1] 2,4-diamino-3,5-diethyltoluene; 2,4-diethyl-6-methyl-1,3-benzenediamine; [2] diethylmethylbenzenediamine [3]	218-255-3 [1] 218-256-9 [2] 270-877-4 [3]	2095-01-4 [1] 2095-02-5 [2] 68479-98-1 [3]	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H373 ** H319 H400 H410	GHS08 GHS07 GHS09 Wng	H312 H302 H373 ** H319 H410			C
612-131-00-6	didecyldimethylammonium chloride	230-525-2	7173-51-5	Acute Tox. 4 * Skin Corr. 1B	H302 H314	GHS05 GHS07 Dgr	H302 H314			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-132-00-1	*N*,*N*'-diphenyl-*p*-phenylenediamine; *N*,*N*'-diphenyl-1,4-benzenediamine	200-806-4	74-31-7	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
612-133-00-7	(4-ammonio-*m*-tolyl)ethyl(2-hydroxyethyl) ammonium sulphate; 4-(*N*-ethyl-*N*-2-hydroxyethyl)-2-methylphenylenediamine sulphate	247-162-0	25646-77-9	Acute Tox. 3 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H301 H373 ** H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H301 H373 ** H317 H410			
612-134-00-2	*N*-(2-(4-amino-*N*-ethyl-*m*-toluidino)ethyl) methanesulphonamide sesquisulphate; 4-(*N*-ethyl-*N*-2-methanesulphonylaminoethyl)-2-methylphenylenediamine sesquisulphate monohydrate	247-161-5	25646-71-3	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
612-135-00-8	*N*-2-naphthylaniline; *N*-phenyl-2-naphthylamine	205-223-9	135-88-6	Carc. 2 Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H351 H319 H315 H317 H411	GHS08 GHS07 GHS09 Wng	H351 H319 H315 H317 H411			
612-136-00-3	*N*-isopropyl-*N*'-phenyl-*p*-phenylenediamine	202-969-7	101-72-4	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410		Skin Sens. 1; H317: C ≥ 0,1 %	
612-137-00-9	4-chloroaniline	203-401-0	106-47-8	Carc. 1B Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H331 H311 H301 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H350 H331 H311 H301 H317 H410			

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612-138-00-4	furalaxyl (ISO); methyl *N*-(2,6-dimethylphenyl)-*N*-(2-furylcarbonyl)-DL-alaninate	260-875-1	57646-30-7	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
612-139-00-X	mefenacet (ISO); 2-(benzothiazol-2-yloxy)-*N*-methyl-*N*-phenylacetamide	277-328-8	73250-68-7	Aquatic Chronic 2	H411	GHS09	H411			
612-140-00-5	quaternary ammonium compounds, benzyl-C_{8-18}-alkyldimethyl, chlorides	264-151-6	63449-41-2	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Aquatic Acute 1	H312 H302 H314 H400	GHS05 GHS07 GHS09 Dgr	H312 H302 H314 H400			
612-141-00-0	4,4'-methylenebis(2-ethylaniline); 4,4'-methylenebis(2-ethylbenzeneamine)	243-420-1	19900-65-3	Carc. 2 Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H351 H302 H400 H410	GHS08 GHS07 GHS09 Wng	H351 H302 H410			
612-142-00-6	biphenyl-2-ylamine	201-990-9	90-41-5	Carc. 2 Acute Tox. 4 * Aquatic Chronic 3	H351 H302 H412	GHS08 GHS07 Wng	H351 H302 H412			
612-143-00-1	N^5,N^5-diethyltoluene-2,5-diamine monohydrochloride; 4-diethylamino-2-methylaniline monohydrochloride	218-130-3	2051-79-8	Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H301 H319 H317 H400 H410	GHS06 GHS09 Dgr	H301 H319 H317 H410			
612-144-00-7	flumetralin (ISO); *N*-(2-chloro-6-fluorobenzyl)-*N*-ethyl-α,α,α-trifluoro-2,6-dinitro-*p*-toluidine	—	62924-70-3	Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H319 H315 H317 H400 H410	GHS07 GHS09 Wng	H319 H315 H317 H410			

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612-145-00-2	*o*-phenylenediamine	202-430-6	95-54-5	Carc. 2 Muta. 2 Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H341 H301 H332 H312 H319 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H351 H341 H301 H332 H312 H319 H317 H410			
612-146-00-8	*o*-phenylenediamine dihydrochloride	210-418-7	615-28-1	Carc. 2 Muta. 2 Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H341 H301 H332 H312 H319 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H351 H341 H301 H332 H312 H319 H317 H410			
612-147-00-3	*m*-phenylenediamine	203-584-7	108-45-2	Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H341 H331 H311 H301 H319 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H341 H331 H311 H301 H319 H317 H410			

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612-148-00-9	*m*-phenylenediamine dihydrochloride	208-790-0	541-69-5	Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H341 H331 H311 H301 H319 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H341 H331 H311 H301 H319 H317 H410			
612-149-00-4	1,3-diphenylguanidine	203-002-1	102-06-7	Repr. 2 Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Chronic 2	H361f *** H302 H319 H335 H315 H411	GHS08 GHS07 GHS09 Wng	H361f *** H302 H319 H335 H315 H411			
612-150-00-X	spiroxamine (ISO); 8-*tert*-butyl-1,4-dioxaspiro[4.5]decan-2-ylmethyl(ethyl)(propyl)amine	—	118134-30-8	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H312 H302 H315 H317 H400 H410	GHS07 GHS09 Wng	H332 H312 H302 H315 H317 H410			
612-151-00-5	diaminotoluene, technical product - reaction mass of [2] and [3]; methyl-phenylenediamine; [1] 4-methyl-*m*-phenylene diamine; [2] 2-methyl-*m*-phenylene diamine [3]	246-910-3 [1] 202-453-1 [2] 212-513-9 [3]	25376-45-8 [1] 95-80-7 [2] 823-40-5 [3]	Carc. 1B Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H350 H301 H332 H312 H319 H317 H411	GHS06 GHS08 GHS09 Dgr	H350 H301 H332 H312 H319 H317 H411			

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612-152-00-0	*N*,*N*-diethyl-*N*',*N*'-dimethylpropan-1,3-diyldiamine	406-610-7	62478-82-4	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1A Aquatic Chronic 3	H226 H332 H302 H373 ** H314 H412	GHS02 GHS08 GHS05 GHS07 Dgr	H226 H332 H302 H373 ** H314 H412			
612-153-00-6	4-[*N*-ethyl-*N*-(2-hydroxyethyl)amino]-1-(2-hydroxyethyl)amino-2-nitrobenzene, monohydrochloride	407-020-2	132885-85-9	Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 3	H302 H317 H412	GHS07 Wng	H302 H317 H412			
612-154-00-1	6'-(isobutylethylamino)-3'-methyl-2'-phenylamino-spiro[isobenzo-2-oxofuran-7,9'-[9*H*]-xanthene]	410-890-6	95235-29-3	Aquatic Chronic 4	H413	—	H413			
612-155-00-7	2'-anilino-6'-((3-ethoxypropyl)ethylamino)-3'-methylspiro(isobenzo-3-oxofuran)-1-(1*H*)-9'-xanthene	411-730-8	93071-94-4	Aquatic Chronic 4	H413	—	H413			
612-156-00-2	reaction mass of: trihexadecylmethylammonium chloride; dihexadecyldimethylammonium chloride	405-620-9	—	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
612-157-00-8	(Z)-1-benzo[*b*]thien-2-ylethanone oxime hydrochloride	410-780-8	—	Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H373 ** H318 H317 H411	GHS08 GHS05 GHS07 GHS09 Dgr	H302 H373 ** H318 H317 H411			
612-158-00-3	reaction mass of: bis(5-dodecyl-2-hydroxybenzald-oximate) copper (II) C_{12}-alkyl group is branched; 4-dodecylsalicylaldoxime	410-820-4	—	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-159-00-9	Reaction products of: trimethylhexamethylene diamine (a mixture of 2,2,4-trimethyl-1,6-hexanediamine and 2,4,4-trimethyl-1,6-hexanediamine, EINECS listed), Epoxide 8 (mono[(C_{10}-C_{16}-alkyloxy)methyl]oxirane derivatives) and *p*-toluene-sulfonic acid	410-880-1	—	Acute Tox. 4 * Skin Corr. 1B Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H314 H410			
612-160-00-4	*p*-toluidine; 4-aminotoluene; [1] toluidinium chloride; [2] toluidine sulphate (1:1) [3]	203-403-1 [1] 208-740-8 [2] 208-741-3 [3]	106-49-0 [1] 540-23-8 [2] 540-25-0 [3]	Carc. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1	H351 H331 H311 H301 H319 H317 H400	GHS06 GHS08 GHS09 Dgr	H351 H331 H311 H301 H319 H317 H400			
612-161-00-X	2,6-xylidine; 2,6-dimethylaniline	201-758-7	87-62-7	Carc. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * STOT SE 3 Skin Irrit. 2 Aquatic Chronic 2	H351 H332 H312 H302 H335 H315 H411	GHS08 GHS07 GHS09 Wng	H351 H332 H312 H302 H335 H315 H411			
612-162-00-5	dimethyldioctadecylammonium chloride; DODMAC	203-508-2	107-64-2	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
612-163-00-0	metalaxyl-M (ISO); mefenoxam; (*R*)-2-[(2,6-dimethylphenyl)-methoxyacetylamino]propionic acid methyl ester	—	70630-17-0	Acute Tox. 4 * Eye Dam. 1	H302 H318	GHS05 GHS07 Dgr	H302 H318			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-164-00-6	2-butyl-2-ethyl-1,5-diaminopentane	412-700-7	137605-95-9	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 3	H312 H302 H373 ** H314 H317 H412	GHS08 GHS05 GHS07 Dgr	H312 H302 H373 ** H314 H317 H412			
612-165-00-1	*N,N'*-diphenyl-*N,N'*-bis(3-methylphenyl)-(1,1'-diphenyl)-4,4'-diamine	413-810-8	65181-78-4	Aquatic Chronic 2	H411	GHS09	H411			
612-166-00-7	reaction mass of: *cis*-(5-ammonium-1,3,3-trimethyl)-cyclohexanemethylammonium phosphate (1:1); *trans*-(5-ammonium-1,3,3-trimethyl)-cyclohexanemethylammonium phosphate (1:1)	411-830-1	114765-88-7	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H318 H317 H412	GHS05 GHS07 Dgr	H318 H317 H412			
612-167-00-2	5-acetyl-3-amino-10,11-dihydro-5*H*-dibenz[*b,f*]azepine-hydrochloride	410-490-1	—	Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H373 ** H318 H317 H411	GHS08 GHS05 GHS07 GHS09 Dgr	H302 H373 ** H318 H317 H411			
612-168-00-8	3,5-dichloro-2,6-difluoropyrdine-4-amine	220-630-1	2840-00-8	Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 2	H312 H302 H411	GHS07 GHS09 Wng	H312 H302 H411			
612-170-00-9	4-chlorophenyl cyclopropyl ketone *O*-(4-aminobenzyl)oxime	405-260-2	—	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
612-171-00-4	*N,N,N',N'*-tetraglycidyl-4,4'-diamino-3,3'-diethyldiphenylmethane	410-060-3	130728-76-6	Muta. 2 Skin Sens. 1 Aquatic Chronic 2	H341 H317 H411	GHS08 GHS09 Wng	H341 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-172-00-X	4,4'-methylenebis(*N*,*N*'-dimethylcyclohexanamine	412-840-9	13474-64-1	Acute Tox. 4 * STOT RE 2 * Skin Corr. 1A Aquatic Chronic 3	H302 H373 ** H314 H412	GHS08 GHS05 GHS07 Dgr	H302 H373 ** H314 H412			
612-173-00-5	lithium 1-amino-4-(4-*tert*-butylanilino) anthraquinone-2-sulfonate	411-140-0	125328-86-1	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			
612-174-00-0	4,4-dimethoxybutylamine	407-690-6	19060-15-2	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 3	H302 H314 H317 H412	GHS05 GHS07 Dgr	H302 H314 H317 H412			
612-175-00-6	2-(*O*-aminooxy)ethylamine dihydrochloride	412-310-7	37866-45-8	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
612-176-00-1	polymer of 1,3-dibromopropane and *N*,*N*-diethyl-*N*',*N*'-dimethyl-1,3-propanediamine	410-570-6	143747-73-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
612-177-00-7	2-naphthylamino-6-sulfomethylamide	412-120-4	104295-55-8	STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H373 ** H317 H411	GHS08 GHS09 Wng	H373 ** H317 H411			
612-178-00-2	1,4,7,10-tetraazacyclododecane disulfate	412-080-8	112193-77-8	Acute Tox. 4 * STOT SE 3 Eye Dam. 1 Aquatic Chronic 3	H302 H335 H318 H412	GHS05 GHS07 Dgr	H302 H335 H318 H412			
612-179-00-8	1-(2-propenyl)pyridinium chloride	412-740-5	25965-81-5	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-180-00-3	3-aminobenzylamine	412-230-2	4403-70-7	Acute Tox. 4 * Skin Corr. 1B Aquatic Chronic 2	H302 H314 H411	GHS05 GHS07 GHS09 Dgr	H302 H314 H411			
612-181-00-9	2-phenylthioaniline	413-030-8	1134-94-7	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
612-182-00-4	1-ethyl-1-methylmorpholinium bromide	418-210-1	65756-41-4	Muta. 2	H341	GHS08 Wng	H341			
612-183-00-X	1-ethyl-1-methylpyrrolidinium bromide	418-200-5	69227-51-6	Muta. 2	H341	GHS08 Wng	H341			
612-184-00-5	6'-(dibutylamino)-3'-methyl-2'-(phenylamino)spiro[isobenzofuran-1(3*H*),9-(9*H*)-xanthen]-3-one	403-830-5	89331-94-2	Aquatic Chronic 3	H412	—	H412			
612-185-00-0	1-[3-[4-((heptadecafluorononyl)oxy)-benzamido]propyl]-*N*,*N*,*N*-trimethylammonium iodide	407-400-8	59493-72-0	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			
612-186-00-6	bis(*N*-(7-hydroxy-8-methyl-5-phenylphenazin-3-ylidene)dimethylammonium) sulfate	406-770-8	149057-64-7	STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H318 H317 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H373 ** H318 H317 H410			
612-187-00-1	2,3,4-trifluoroaniline	407-170-9	3862-73-5	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Irrit. 2 Eye Dam. 1 Aquatic Chronic 2	H312 H302 H373 ** H315 H318 H411	GHS08 GHS05 GHS07 GHS09 Dgr	H312 H302 H373 ** H315 H318 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-188-00-7	4,4'-(9*H*-fluoren-9-ylidene)bis(2-chloroaniline)	407-560-9	107934-68-9	Aquatic Chronic 2	H411	GHS09	H411			
612-189-00-2	4-amino-2-(aminomethyl)phenol dihydrochloride	412-510-4	135043-64-0	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
612-190-00-8	4,4'-methylenebis(2-isopropyl-6-methylaniline)	415-150-6	16298-38-7	STOT RE 2 * Aquatic Chronic 2	H373 ** H411	GHS08 GHS09 Wng	H373 ** H411			
612-191-00-3	polymer of allylamine hydrochloride	415-050-2	71550-12-4	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			
612-192-00-9	2-isopropyl-4-(*N*-methyl)aminomethylthiazole	414-800-6	154212-60-9	Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Eye Dam. 1 Aquatic Chronic 2	H312 H302 H315 H318 H411	GHS05 GHS07 GHS09 Dgr	H312 H302 H315 H318 H411			
612-193-00-4	3-methylaminomethylphenylamine	414-570-7	18759-96-1	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H312 H302 H314 H317 H410			
612-194-00-X	2-hydroxy-3-[(2-hydroxyethyl)-[2-(1-oxotetradecyl)amino]ethyl]amino]-*N*,*N*,*N*-trimethyl-1-propanammonium chloride	414-670-0	141890-30-4	Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H318 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H318 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-195-00-5	bis[tributyl 4-(methylbenzyl)ammonium] 1,5-naphthalenedisulfonate	415-210-1	160236-81-7	Acute Tox. 4 * Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H302 H318 H400 H410	GHS05 GHS07 GHS09 Dgr	H332 H302 H318 H410			
612-196-00-0	4-chloro-*o*-toluidine; [1] 4-chloro-*o*-toluidine hydrochloride [2]	202-441-6 [1] 221-627-8 [2]	95-69-2 [1] 3165-93-3 [2]	Carc. 1B Muta. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H350 H341 H331 H311 H301 H400 H410	GHS06 GHS08 GHS09 Dgr	H350 H341 H331 H311 H301 H410			
612-197-00-6	2,4,5-trimethylaniline; [1] 2,4,5-trimethylaniline hydrochloride [2]	205-282-0 [1] - [2]	137-17-7 [1] 21436-97-5 [2]	Carc. 1B Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H350 H331 H311 H301 H411	GHS06 GHS08 GHS09 Dgr	H350 H331 H311 H301 H411			
612-198-00-1	4,4'-thiodianiline and its salts	205-370-9	139-65-1	Carc. 1B Acute Tox. 4 * Aquatic Chronic 2	H350 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H302 H411			
612-199-00-7	4,4'-oxydianiline and its salts; *p*-aminophenyl ether	202-977-0	101-80-4	Carc. 1B Muta. 1B Repr. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 2	H350 H340 H361f *** H331 H311 H301 H411	GHS06 GHS08 GHS09 Dgr	H350 H340 H361f *** H331 H311 H301 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-200-00-0	2,4-diaminoanisole; 4-methoxy-*m*-phenylenediamine; [1] 2,4-diaminoanisole sulphate [2]	210-406-1 [1] 254-323-9 [2]	615-05-4 [1] 39156-41-7 [2]	Carc. 1B Muta. 2 Acute Tox. 4 * Aquatic Chronic 2	H350 H341 H302 H411	GHS08 GHS07 GHS09 Dgr	H350 H341 H302 H411			
612-201-00-6	N,N,N',N'-tetramethyl-4,4'-methylendianiline	202-959-2	101-61-1	Carc. 1B Aquatic Acute 1 Aquatic Chronic 1	H350 H400 H410	GHS08 GHS09 Dgr	H350 H410			
612-202-00-1	3,4-dichloroaniline	202-448-4	95-76-1	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H318 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H331 H311 H301 H318 H317 H410			
612-204-00-2	C.I. Basic Violet 3; 4-[4,4'-bis(dimethylamino) benzhydrylidene]cyclohexa-2,5-dien-1-ylidene]dimethylammonium chloride	208-953-6	548-62-9	Carc. 2 Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H302 H318 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H351 H302 H318 H410			
612-205-00-8	C.I. Basic Violet 3 with ≥ 0,1 % of Michler's ketone (EC no. 202-027-5)	208-953-6	548-62-9	Carc. 1B Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H302 H318 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H350 H302 H318 H410			
612-206-00-3	famoxadone (ISO); 3-anilino-5-methyl-5-(4-phenoxyphenyl)-1,3-oxazolidine-2,4-dione	—	131807-57-3	STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H373 ** H400 H410	GHS08 GHS09 Wng	H373 ** H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-207-00-9	4-ethoxyaniline; *p*-phenetidine	205-855-5	156-43-4	Muta. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1	H341 H332 H312 H302 H319 H317	GHS08 GHS07 Wng	H341 H332 H312 H302 H319 H317			
612-209-00-X	6-methoxy-*m*-toluidine; *p*-cresidine	204-419-1	120-71-8	Carc. 1B Acute Tox. 4 *	H350 H302	GHS08 GHS07 Dgr	H350 H302			
612-210-00-5	5-nitro-*o*-toluidine; [1] 5-nitro-*o*-toluidine hydrochloride [2]	202-765-8 [1] 256-960-8 [2]	99-55-8 [1] 51085-52-0 [2]	Carc. 2 Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Aquatic Chronic 3	H351 H331 H311 H301 H412	GHS06 GHS08 Dgr	H351 H331 H311 H301 H412			
612-211-00-0	N-[(benzotriazole-1-yl)methyl)]-4-carboxybenzenesulfonamide	416-470-9	170292-97-4	Eye Irrit. 2 Aquatic Chronic 2	H319 H411	GHS07 GHS09 Wng	H319 H411			
612-212-00-6	2,6-dichloro-4-trifluoromethylaniline	416-430-0	24279-39-8	Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H302 H315 H317 H400 H410	GHS07 GHS09 Wng	H332 H302 H315 H317 H410			
612-213-00-1	isobutylidene-(2-(2-isopropyl-4,4-dimethyloxazolidine-3-yl)-1,1-dimethylethyl)amine	419-850-2	148348-13-4	Skin Corr. 1B Aquatic Chronic 3	H314 H412	GHS05 Dgr	H314 H412			
612-214-00-7	4-(2,2-diphenylethenyl)-*N*,*N*-di-phenylbenzenamine	421-390-2	89114-90-9	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
612-215-00-2	3-chloro-2-(isopropylthio)aniline	421-700-6	179104-32-6	Skin Irrit. 2 Aquatic Chronic 2	H315 H411	GHS07 GHS09 Wng	H315 H411			
612-217-00-3	1-methoxy-2-propylamine	422-550-4	37143-54-7	Flam. Liq. 2 Skin Corr. 1B Acute Tox. 4 * Aquatic Chronic 3	H225 H314 H302 H412	GHS02 GHS05 GHS07 Dgr	H225 H314 H302 H412			
613-001-00-1	ethyleneimine; aziridine	205-793-9	151-56-4	Flam. Liq. 2 Carc. 1B Muta. 1B Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Skin Corr. 1B Aquatic Chronic 2	H225 H350 H340 H330 H310 H300 H314 H411	GHS02 GHS06 GHS08 GHS05 GHS09 Dgr	H225 H350 H340 H330 H310 H300 H314 H411			D
613-002-00-7	pyridine	203-809-9	110-86-1	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H225 H332 H312 H302	GHS02 GHS07 Dgr	H225 H332 H312 H302		*	
613-003-00-2	1,2,3,4-tetranitrocarbazole	—	6202-15-9	Expl. 1.1 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 *	H201 H332 H312 H302	GHS01 GHS07 Wng	H201 H332 H312 H302			
613-004-00-8	crimidine (ISO); 2-chloro-6-methylpyrimidin-4-yldimethylamine	208-622-6	535-89-7	Acute Tox. 2 *	H300	GHS06 Dgr	H300			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-007-00-4	desmetryne (ISO); 6-isopropylamino-2-methylamino-4-methylthio-1,3,5-triazine	213-800-1	1014-69-3	Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H400 H410	GHS07 GHS09 Wng	H312 H302 H410			
613-008-00-X	dazomet (ISO); tetrahydro-3,5-dimethyl-1,3,5-thiadiazine-2-thione	208-576-7	533-74-4	Acute Tox. 4 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H302 H319 H400 H410	GHS07 GHS09 Wng	H302 H319 H410			
613-009-00-5	2,4,6-trichloro-1,3,5-triazine; cyanuric chloride	203-614-9	108-77-0	Acute Tox. 2 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1	H330 H302 H314 H317	GHS06 GHS05 Dgr	H330 H302 H314 H317	EUH014	STOT SE 3; H335: C ≥ 5 %	
613-010-00-0	ametryn (ISO); 2-ethylamino-4-isopropylamino-6-methylthio-1,3,5-triazine	212-634-7	834-12-8	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-011-00-6	amitrole (ISO); 1,2,4-triazol-3-ylamine	200-521-5	61-82-5	Repr. 2 STOT RE 2 * Aquatic Chronic 2	H361d *** H373 ** H411	GHS08 GHS09 Wng	H361d *** H373 ** H411			
613-012-00-1	bentazone (ISO); 3-isopropyl-2,1,3-benzothiadiazine-4-one-2,2-dioxide	246-585-8	25057-89-0	Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 3	H302 H319 H317 H412	GHS07 Wng	H302 H319 H317 H412			
613-013-00-7	cyanazine (ISO); 2-(4-chloro-6-ethylamino-1,3,5-triazine-2-ylamino)-2-methylpropionitrile	244-544-9	21725-46-2	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-014-00-2	ethoxyquin(ISO); 6-ethoxy-1,2-dihydro-2,2,4-trimethylquinoline	202-075-7	91-53-2	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-015-00-8	fenazaflor (ISO); phenyl 5,6-dichloro-2-trifluoromethylbenzimidazole-1-carboxylate	238-134-9	14255-88-0	Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H400 H410	GHS07 GHS09 Wng	H312 H302 H410			
613-016-00-3	fuberidazole(ISO); 2-(2-furyl)benzimidazole	223-404-0	3878-19-1	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-017-00-9	bis (8-hydroxyquinolinium) sulphate	205-137-1	134-31-6	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-018-00-4	morfamquat (ISO); 1,1'-bis(3,5-dimethylmorpholinocarbonylmethyl)-4,4'-bipyridilium ion	—	7411-47-4	Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Chronic 3	H302 H319 H335 H315 H412	GHS07 Wng	H302 H319 H335 H315 H412			
613-019-00-X	thioquinox(ISO); 2-thio-1,3-dithiolo(4,5,b)quinoxaline	202-272-8	93-75-4	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-020-00-5	tridemorph (ISO); 2,6-dimethyl-4-tridecylmorpholine	246-347-3	24602-86-6	Repr. 1B Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H332 H302 H315 H400 H410	GHS08 GHS07 GHS09 Dgr	H360D *** H332 H302 H315 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-021-00-0	dithianon (ISO); 5,10-dihydro-5,10-dioxonaphtho(2,3-b)(1,4)dithiazine-2,3-dicarbonitrile	222-098-6	3347-22-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-022-00-6	pyrethrins including cinerins, with the exception of those specified elsewhere in this Annex	—	—	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H332 H312 H302 H400 H410	GHS07 GHS09 Wng	H332 H312 H302 H410			A
613-023-00-1	2-methyl-4-oxo-3-(penta-2,4-dienyl)cyclopent-2-enyl [1*R*-[1α[*S**(*Z*)],3β]]-chrysanthemate; pyrethrin I	204-455-8	121-21-1	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H332 H312 H302 H400 H410	GHS07 GHS09 Wng	H332 H312 H302 H410			
613-024-00-7	2-methyl-4-oxo-3-(penta-2,4-dienyl)cyclopent-2-enyl[1*R*-[1α[*S**(*Z*)](3β)]]-3-(3-methoxy-2-methyl-3-oxoprop-1-enyl)-2,2-dimethylcyclopropanecarboxylate; pyrethrin II	204-462-6	121-29-9	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H332 H312 H302 H400 H410	GHS07 GHS09 Wng	H332 H312 H302 H410			
613-025-00-2	cinerin I; 3-(but-2-enyl)-2-methyl-4-oxocyclopent-2-enyl 2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate	246-948-0	25402-06-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-026-00-8	cinerin II; 3-(but-2-enyl)-2-methyl-4-oxocyclopent-2-enyl 2,2-dimethyl-3-(3-methoxy-2-methyl-3-oxoprop-1-enyl)cyclopropanecarboxylate	204-454-2	121-20-0	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-027-00-3	piperidine	203-813-0	110-89-4	Flam. Liq. 2 Acute Tox. 3 * Acute Tox. 3 * Skin Corr. 1B	H225 H331 H311 H314	GHS02 GHS06 GHS05 Dgr	H225 H331 H311 H314		*	
613-028-00-9	morpholine	203-815-1	110-91-8	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H226 H332 H312 H302 H314	GHS02 GHS05 GHS07 Dg	H226 H332 H312 H302 H314			
613-029-00-4	dichloro-1,3,5-triazinetrione; dichloroisocyanuric acid	220-487-5	2782-57-2	Ox. Sol. 2 Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Aquatic Acute 1 Aquatic Chronic 1	H272 H302 H319 H335 H400 H410	GHS03 GHS07 GHS09 Dgr	H272 H302 H319 H335 H410	EUH031		T
613-030-00-X	troclosene potassium; [1] troclosene sodium [2]	218-828-8 [1] 220-767-7 [2]	2244-21-5 [1] 2893-78-9 [2]	Ox. Sol. 2 Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Aquatic Acute 1 Aquatic Chronic 1	H272 H302 H319 H335 H400 H410	GHS03 GHS07 GHS09 Dgr	H272 H302 H319 H335 H410	EUH031	* STOT SE 3; H335: C ≥ 10 % EUH031: C ≥ 10 %	
613-030-01-7	troclosene sodium, dihydrate	220-767-7	51580-86-0	Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Aquatic Acute 1 Aquatic Chronic 1	H302 H319 H335 H400 H410	GHS07 GHS09 Wng	H302 H319 H335 H410	EUH031		

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-031-00-5	symclosene; trichloroisocyanuric acid; trichloro-1,3,5-triazinetrion	201-782-8	87-90-1	Ox. Sol. 2 Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Aquatic Acute 1 Aquatic Chronic 1	H272 H302 H319 H335 H400 H410	GHS03 GHS07 GHS09 Dgr	H272 H302 H319 H335 H410	EUH031		
613-032-00-0	methyl-2,3,5,6-tetrachloro-4-pyridylsulphone; 2,3,5,6-tetrachloro-4-(methylsulphonyl)pyridine	236-035-5	13108-52-6	Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1	H312 H302 H319 H317	GHS07 Wng	H312 H302 H319 H317			
613-033-00-6	2-methylaziridine; propyleneimine	200-878-7	75-55-8	Flam. Liq. 2 Carc. 1B Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Eye Dam. 1 Aquatic Chronic 2	H225 H350 H330 H310 H300 H318 H411	GHS02 GHS06 GHS08 GHS05 GHS09 Dgr	H225 H350 H330 H310 H300 H318 H411		Carc. 1B; H350: C ≥ 0,01 %	
613-034-00-1	1,2-dimethylimidazole	217-101-2	1739-84-0	Acute Tox. 4 * Skin Irrit. 2 Eye Dam. 1	H302 H315 H318	GHS05 GHS07 Dgr	H302 H315 H318			
613-035-00-7	1-methylimidazole	210-484-7	616-47-7	Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1B	H312 H302 H314	GHS05 GHS07 Dgr	H312 H302 H314			
613-036-00-2	2-methylpyridine; 2-picoline	203-643-7	109-06-8	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 STOT SE 3	H226 H332 H312 H302 H319 H335	GHS02 GHS07 Wng	H226 H332 H312 H302 H319 H335			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-037-00-8	4-methylpyridine; 4-picoline	203-626-4	108-89-4	Flam. Liq. 3 Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2	H226 H311 H332 H302 H319 H335 H315	GHS02 GHS06 Dgr	H226 H311 H332 H302 H319 H335 H315			
613-038-00-3	6-phenyl-1,3,5-triazine-2,4-diyldiamine; 6-phenyl-1,3,5-triazine-2,4-diamine; benzoguanamine	202-095-6	91-76-9	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
613-039-00-9	ethylene thiourea; imidazolidine-2-thione; 2-imidazoline-2-thiol	202-506-9	96-45-7	Repr. 1B Acute Tox. 4 *	H360-D *** H302	GHS08 GHS07 Dgr	H360D *** H302			
613-040-00-4	azaconazole (ISO); 1-{[2-(2,4-dichlorophenyl)-1,3-dioxolan-2-yl]methyl}-1*H*-1,2.4-triazole	262-102-3	60207-31-0	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-041-00-X	morpholine-4-carbonyl chloride	239-213-0	15159-40-7	Carc. 2 Eye Irrit. 2 Skin Irrit. 2	H351 H319 H315	GHS08 Wng	H351 H319 H315	EUH014		
613-042-00-5	imazalil (ISO); 1-[2-(allyloxy)-2-(2,4-dichlorophenyl)ethyl]-1*H*-imidazole	252-615-0	35554-44-0	Acute Tox. 4 * Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H302 H318 H400 H410	GHS05 GHS07 GHS09 Dgr	H332 H302 H318 H410			
613-043-00-0	imazalil sulphate (ISO) powder; 1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate; [1] (±)-1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1H-imidazolium hydrogen sulphate [2]	261-351-5 [1] 281-291-3 [2]	58594-72-2 [1] 83918-57-4 [2]	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-043-01-8	imazalil sulphate (ISO), aqueous solution; 1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate; [1] (±)-1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate [2]	261-351-5 [1] 281-291-3 [2]	58594-72-2 [1] 83918-57-4 [2]	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Wng	H302 H314 H317 H410		Skin Corr. 1B; H314: C ≥ 50 % Skin Irrit. 2; H315: 30 % ≤ C < 50 % Eye Dam. 1; H318: 15 % ≤ C < 50 % Eye Irrit. 2; H319: 5 % ≤ C < 15 %	
613-044-00-6	captan (ISO); 1,2,3,6-tetrahydro-*N*-(trichloromethylthio) phthalimide	205-087-0	133-06-2	Carc. 2 Acute Tox. 3 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H351 H331 H318 H317 H400	GHS06 GHS08 GHS05 GHS09 Dgr	H351 H331 H318 H317 H400			
613-045-00-1	folpet (ISO); *N*-(trichloromethylthio)phthalimide	205-088-6	133-07-3	Carc. 2 Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1	H351 H332 H319 H317 H400	GHS08 GHS07 GHS09 Wng	H351 H332 H319 H317 H400			
613-046-00-7	captafol (ISO); 1,2,3,6-tetrahydro-*N*-(1,1,2,2-tetrachloroethylthio)phthalimide	219-363-3	2425-06-1	Carc. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H350 H317 H400 H410	GHS08 GHS09 Dgr	H350 H317 H410			
613-047-00-2	1-dimethylcarbamoyl-5-methylpyrazol-3-yl dimethylcarbamate; dimetilan (ISO)	211-420-0	644-64-4	Acute Tox. 3 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H301 H312 H400 H410	GHS06 GHS09 Dgr	H301 H312 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-048-00-8	carbendazim (ISO); methyl benzimidazol-2-ylcarbamate	234-232-0	10605-21-7	Muta. 1B Repr. 1B Aquatic Acute 1 Aquatic Chronic 1	H340 H360-FD H400 H410	GHS08 GHS09 Dgr	H340 H360FD H410			
613-049-00-3	benomyl (ISO); methyl 1-(butylcarbamoyl)benzimidazol-2-ylcarbamate	241-775-7	17804-35-2	Muta. 1B Repr. 1B STOT SE 3 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H340 H360-FD H335 H315 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H340 H360FD H335 H315 H317 H410		M = 10	
613-050-00-9	carbadox (INN); methyl 3-(quinoxalin-2-ylmethylene)carbazate 1,4-dioxide; 2-(methoxycarbonylhydrazonomethyl)quinoxaline 1,4-dioxide	229-879-0	6804-07-5	Flam. Sol. 1 Carc. 1B Acute Tox. 4 *	H228 H350 H302	GHS02 GHS08 GHS07 Dgr	H228 H350 H302			T
613-051-00-4	molinate (ISO); *S*-ethyl 1-perhydroazepinecarbothioate; *S*-ethyl perhydroazepine-1-carbothioate	218-661-0	2212-67-1	Carc. 2 Repr. 2 Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H361f *** H332 H302 H373 ** H317 H400 H410	GHS08 GHS07 GHS09 Wng	H351 H361f *** H332 H302 H373 ** H317 H410		M = 100	
613-052-00-X	trifenmorph (ISO); 4-tritylmorpholine	215-812-2	1420-06-0	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-053-00-5	anilazine (ISO); 2-chloro-*N*-(4,6-dichloro-1,3,5-triazin-2-yl) aniline	202-910-5	101-05-3	Eye Irrit. 2 Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H319 H315 H400 H410	GHS07 GHS09 Wng	H319 H315 H410			
613-054-00-0	thiabendazol (ISO); 2-(thiazole-4-yl)benzimidazole	205-725-8	148-79-8	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-056-00-1	1,2-dimethyl-3,5-diphenylpyrazolium methylsulphate; difenzoquat methyl sulfate	256-152-5	43222-48-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS09 Wng	H302 H410			
613-057-00-7	dodemorph (ISO); 4-cyclododecyl-2,6-dimethylmorpholine	216-474-9	1593-77-7	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Chronic 2	H319 H335 H315 H411	GHS07 GHS09 Wng	H319 H335 H315 H411			
613-058-00-2	permethrin (ISO); *m*-phenoxybenzyl 3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate	258-067-9	52645-53-1	Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H302 H317 H400 H410	GHS07 GHS09 Wng	H332 H302 H317 H410		M = 1 000	
613-059-00-8	profluralin (ISO); *N*-(cyclopropylmethyl)-α,α,α-trifluoro-2,6-dinitro-*N*-propyl-*p*-toluidine	247-656-6	26399-36-0	Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H319 H400 H410	GHS07 GHS09 Wng	H319 H410			
613-060-00-3	resmethrin (ISO); 5-benzyl-3-furylmethyl (±)-*cis-trans*-chrysanthemate	233-940-7	10453-86-8	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-061-00-9	6-(1α,5aβ,8aβ,9-pentahydroxy-7β-isopropyl-2β,5β,8β-trimethylperhydro-8bα,9-epoxy-5,8-ethanocyclopenta[1,2-b]indenyl) pyrrole-2-carboxylate; ryania	239-732-2	15662-33-6	Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H400 H410	GHS07 GHS09 Wng	H312 H302 H410			
613-062-00-4	sabadilla (ISO); veratrine	—	8051-02-3	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2	H319 H335 H315	GHS07 Wng	H319 H335 H315			
613-063-00-X	secbumeton (ISO); 2-sec-butylamino-4-ethylamino-6-methoxy-1,3,5-triazine	247-554-1	26259-45-0	Acute Tox. 4 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H302 H319 H400 H410	GHS07 GHS09 Wng	H302 H319 H410			
613-064-00-5	5-(3,6,9-trioxa-2-undecyloxy)benzo(d)-1,3-dioxolane; sesamex	—	51-14-9	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-065-00-0	simetryn (ISO); 2,4-bis(ethylamino)-6-methylthio-1,3,5-triazine	213-801-7	1014-70-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-066-00-6	terbumeton (ISO); 2-*tert*-butylamino-4-ethylamino-6-methoxy-1,3,5-triazine	251-637-8	33693-04-8	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-067-00-1	propazine(ISO); 2-chloro-4,6-bis(isopropylamino)-1,3,5-triazine	205-359-9	139-40-2	Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H400 H410	GHS08 GHS09 Wng	H351 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-068-00-7	atrazine (ISO); 2-chloro-4-ethylamine-6-isopropylamine-1,3,5-triazine	217-617-8	1912-24-9	STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H317 H400 H410	GHS08 GHS09 Wng	H373 ** H317 H410			
613-069-00-2	ε-caprolactam	203-313-2	105-60-2	Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2	H332 H302 H319 H335 H315	GHS07 Wng	H332 H302 H319 H335 H315			
613-070-00-8	propylenethiourea	—	2122-19-2	Repr. 2 Acute Tox. 4 * Aquatic Chronic 3	H361d *** H302 H412	GHS08 GHS07 Wng	H361d *** H302 H412			
613-071-00-3	2-fluoro-5-trifluoromethylpyridine	400-290-2	69045-82-5	Flam. Liq. 3 Skin Sens. 1 Aquatic Chronic 3	H226 H317 H412	GHS02 GHS07 Wng	H226 H317 H412			
613-072-00-9	*N*,*N*-bis(2-ethylhexyl)-((1,2,4-triazol-1-yl) methyl)amine	401-280-0	91273-04-0	Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 2	H314 H317 H411	GHS05 GHS07 GHS09 Dgr	H314 H317 H411			
613-073-00-4	*N*,*N*-dimethyl-2-(3-(4-chlorophenyl)-4,5-dihydropyrazol-1-ylphenylsulphonyl)ethylamine	401-410-6	10357-99-0	STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H373 ** H317 H411	GHS08 GHS09 Wng	H373 ** H317 H411			
613-074-00-X	3-(3-methylpent-3-yl)isoxazol-5-ylamine	401-460-9	82560-06-3	Acute Tox. 3 * Acute Tox. 3 * Eye Dam. 1 Aquatic Chronic 3	H331 H301 H318 H412	GHS06 GHS05 Dgr	H331 H301 H318 H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-075-00-5	1,3-dichloro-5-ethyl-5-methylimidazolidine-2,4-dione	401-570-7	89415-87-2	Ox. Sol. 1 **** Acute Tox. 3 * Skin Corr. 1B Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1	H271 H331 H314 H302 H317 H400	GHS03 GHS06 GHS05 GHS09 Dgr	H271 H331 H314 H302 H317 H400			
613-076-00-0	3-chloro-5-trifluoromethyl-2-pyridylamine	401-670-0	79456-26-1	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
613-077-00-6	reaction mass of 5-heptyl-1,2,4-triazol-3-ylamine and 5-nonyl-1,2,4-triazol-3-ylamine	401-940-8	—	Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 2	H302 H319 H411	GHS07 GHS09 Wng	H302 H319 H411			
613-078-00-1	*N*,*N*,*N*,*N*-tetrakis(4,6-bis(butyl-(*N*-methyl-2,2,6,6-tetramethylpiperidin-4-yl)amino) triazin-2-yl)-4,7-diazadecane-1,10-diamine	401-990-0	106990-43-6	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-079-00-7	4-(1(or 4 or 5 or 6)-methyl-8,9,10-trinorborn-5-en-2-yl)pyridine, reaction mass of isomers	402-520-7	—	Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H315 H317 H400 H410	GHS07 GHS09 Wng	H312 H302 H315 H317 H410			
613-080-00-2	3-(bis(2-ethylhexyl)aminomethyl)benzothiazole-2(3*H*)-thione	402-540-6	105254-85-1	Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H314 H317 H410			
613-081-00-8	1-butyl-2-methylpyridinium bromide	402-680-8	26576-84-1	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-082-00-3	2-methyl-1-pentylpyridinium bromide	402-690-2	—	Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H312 H302 H412	GHS07 Wng	H312 H302 H412			
613-083-00-9	2-(4-(3-(4-chlorophenyl)-2-pyrazolin-1-yl) phenylsulfonyl)ethyldimethylammonium formate	402-120-2	—	Skin Corr. 1B STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H314 H373 ** H317 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H314 H373 ** H317 H410			
613-084-00-4	2-(4-(3-(4-chlorophenyl)-4,5-dihydropyrazolyl)phenylsulphonyl)ethyldimethylammonium hydrogen phosphonate	402-490-5	106359-93-7	Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H319 H400 H410	GHS07 GHS09 Wng	H319 H410			
613-085-00-X	reaction mass of 1,1'-(methylenebis(4,1-phenylene))dipyrrole-2,5-dione and *N*-(4-(4-(2,5-dioxopyrrol-1-yl)benzyl)phenyl)acetamide and 1-(4-(4-(5-oxo-2*H*-2-furylidenamino)benzyl)phenyl)pyrrole-2,5-dione	401-970-1	—	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
613-086-00-5	caffeine	200-362-1	58-08-2	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-087-00-0	tetrahydrothiophene	203-728-9	110-01-0	Flam. Liq. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 3	H225 H332 H312 H302 H319 H315 H412	GHS02 GHS07 Dgr	H225 H332 H312 H302 H319 H315 H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-088-00-6	1,2-benzisothiazol-3(2*H*)-one; 1,2-benzisothiazolin-3-one	220-120-9	2634-33-5	Acute Tox. 4 * Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H302 H315 H318 H317 H400	GHS05 GHS07 GHS09 Dgr	H302 H315 H318 H317 H400		Skin Sens. 1; H317: C ≥ 0,05 %	
613-089-00-1	diquat dibromide; [1] diquat dichloride; [2] 6,7-dihydrodipyrido[1,2-α:2',1'-c]pyrazinediylium dihydroxide [3]	201-579-4 [1] 223-714-6 [2] 301-467-6 [3]	85-00-7 [1] 4032-26-2 [2] 94021-76-8 [3]	Acute Tox. 2 * STOT RE 1 Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H330 H372 ** H302 H319 H335 H315 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H330 H372 ** H302 H319 H335 H315 H317 H410			
613-090-00-7	paraquat dichloride; 1,1-dimethyl-4,4'-bipyridinium dichloride; [1] paraquat dimethylsulfate; 1,1-dimethyl-4,4'-bipyridinium dimethyl sulphate [2]	217-615-7 [1] 218-196-3 [2]	1910-42-5 [1] 2074-50-2 [2]	Acute Tox. 2 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H330 H311 H301 H372 ** H319 H335 H315 H400 H410	GHS06 GHS08 GHS09 Dgr	H330 H311 H301 H372 ** H319 H335 H315 H410			
613-091-00-2	morfamquat dichloride; [1] morfamquat sulfate [2]	225-062-8 [1] [2]	4636-83-3 [1] 29873-36-7 [2]	Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Chronic 3	H302 H319 H335 H315 H412	GHS07 Wng	H302 H319 H335 H315 H412			

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613-092-00-8	1,10-phenanthroline	200-629-2	66-71-7	Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H301 H400 H410	GHS06 GHS09 Dgr	H301 H410			
613-093-00-3	hexasodium 6,13-dichloro-3,10-bis((4-(2,5-disulfonatoanilino)-6-fluoro-1,3,5-triazin-2-ylamino)prop-3-ylamino)-5,12-dioxa-7,14-diazapentacene-4,11-disulfonate	400-050-7	85153-92-0	Resp. Sens. 1 Skin Sens. 1	H334 H317	GHS08 Dgr	H334 H317			
613-094-00-9	4-methoxy-*N*,6-dimethyl-1,3,5-triazin-2-ylamine	401-360-5	5248-39-5	Acute Tox. 4 * STOT RE 2 *	H302 H373 **	GHS08 GHS07 Wng	H302 H373 **			
613-095-00-4	sodium 3-(2*H*-benzotriazol-2-yl)-5-*sec*-butyl-4-hydroxybenzenesulfonate	403-080-9	92484-48-5	Eye Dam. 1	H318	GHS05 Dgr	H318			
613-096-00-X	2-amino-6-ethoxy-4-methylamino-1,3,5-triazine	403-580-7	62096-63-3	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-097-00-5	7-amino-3-((5-carboxymethyl-4-methyl-1,3-thiazol-2-ylthio)methyl)-8-oxo-5-thia-1-azabicyclo(4.2.0)oct-2-ene-2-carboxylic acid	403-690-5	111298-82-9	Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 3	H334 H317 H412	GHS08 Dgr	H334 H317 H412			
613-098-00-0	*N*-(*n*-octyl)-2-pyrrolidone	403-700-8	2687-94-7	Skin Corr. 1B Aquatic Chronic 2	H314 H411	GHS05 GHS09 Dgr	H314 H411			
613-099-00-6	1-dodecyl-2-pyrrolidone	403-730-1	2687-96-9	Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H314 H317 H410			
613-100-00-X	2,9-bis(3-(diethylamino)propylsulfamoyl) quino(2,3-b)acridine-7,14-dione	404-230-6	—	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-101-00-5	*N*—*tert*-pentyl-2-benzothiazolesulfenamide	404-380-2	110799-28-5	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
613-102-00-0	dimethomorph (ISO); 4-(3-(4-chlorophenyl)-3-(3,4-dimethoxyphenyl)acryloyl)morpholine	404-200-2	110488-70-5	Aquatic Chronic 2	H411	GHS09	H411			
613-103-00-6	sodium 5-*n*-butylbenzotriazole	404-450-2	118685-34-0	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 2	H302 H314 H317 H411	GHS05 GHS07 GHS09 Dgr	H302 H314 H317 H411			
613-104-00-1	5-*tert*-butyl-3-isoxazolylamine hydrochloride	404-840-2	—	Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Aquatic Chronic 3	H302 H373 ** H318 H412	GHS08 GHS05 GHS07 Dgr	H302 H373 ** H318 H412			
613-105-00-7	hexakis(tetramethylammonium) 4,4'-vinylenebis((3-sulfonato-4,1-phenylene)imino (6-morpholino-1,3,5-triazine-4,2-diyl) imino)bis(5-hydroxy-6-phenylazonaphthalene-2,7-disulfonate)	405-160-9	124537-30-0	Acute Tox. 3 * Skin Sens. 1 Aquatic Chronic 3	H301 H317 H412	GHS06 Dgr	H301 H317 H412			
613-106-00-2	tetrapotassium 2-(4-(5-(1-(2,5-disulfonatophenyl)-3-ethoxycarbonyl-5-hydroxypyrazol-4-yl)penta-2,4-dienylidene)-3-ethoxycarbonyl-5-oxo-2-pyrazolin-1-yl)benzene-1,4-disulfonate	405-240-3	—	Skin Sens. 1	H317	GHS07 Wng	H317			
613-107-00-8	hexasodium 2,2'-vinylenebis((3-sulfonato-4,1-phenylene)imino(6-(*N*-cyanoethyl-*N*-(2-hydroxypropyl)amino)-1,3,5-triazine-4,2-diyl)imino)dibenzene-1,4-disulfonate	405-280-1	76508-02-6	Eye Irrit. 2	H319	GHS07 Wng	H319			
613-108-00-3	benzothiazole-2-thiol	205-736-8	149-30-4	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			

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613-109-00-9	bis(piperidinothiocarbonyl) disulphide	202-328-1	94-37-1	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Skin Sens. 1	H319 H335 H315 H317	GHS07 Wng	H319 H335 H315 H317			
613-110-00-4	dimepiperate (ISO); *S*-(1-methyl-1-phenylethyl) piperidine-1-carbothioate	262-784-2	61432-55-1	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
613-111-00-X	1,2,4-triazole	206-022-9	288-88-0	Repr. 2 Acute Tox. 4 * Eye Irrit. 2	H361d *** H302 H319	GHS08 GHS07 Wng	H361d *** H302 H319			
613-112-00-5	octhilinone (ISO); 2-octyl-2*H*-isothiazol-3-one	247-761-7	26530-20-1	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H302 H314 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H331 H311 H302 H314 H317 H410		Skin Sens. 1; H317: C ≥ 0,05 %	
613-113-00-0	2-(morpholinothio)benzothiazole	203-052-4	102-77-2	Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H319 H315 H317 H411	GHS07 GHS09 Wng	H319 H315 H317 H411			
613-114-00-6	2,2',2"-(hexahydro-1,3,5-triazine-1,3,5-triyl) triethanol; 1,3,5-tris(2-hydroxyethyl)hexahydro-1,3,5-triazine	225-208-0	4719-04-4	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317		Skin Sens. 1; H317: C ≥ 0,1 %	
613-115-00-1	hymexazol (ISO); 3-hydroxy-5-methylisoxazole	233-000-6	10004-44-1	Acute Tox. 4 * Eye Dam. 1 Aquatic Chronic 3	H302 H318 H412	GHS05 GHS07 Dgr	H302 H318 H412			

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613-116-00-7	tolylfluanid (ISO); dichloro-*N*-[(dimethylamino)sulphonyl]fluoro-*N*-(*p*-tolyl)methanesulphenamide	211-986-9	731-27-1	Acute Tox. 3 * STOT RE 2 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H373 ** H319 H335 H315 H317 H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H373 ** H319 H335 H315 H317 H410			
613-117-00-2	diniconazole (ISO); (*E*)-β-[(2,4-dichlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*—1,2,4-triazol-1-ethanol; (*E*)-(*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*—1,2,4-triazol-1-yl)pent-1-en-3-ol	—	76714-88-0	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-118-00-8	flubenzimine (ISO); *N*-[3-phenyl-4,5-bis[(trifluoromethyl)imino]thiazolidin-2-ylidene]aniline	253-703-1	37893-02-0	Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H319 H400 H410	GHS07 GHS09 Wng	H319 H410			
613-119-00-3	(benzothiazol-2-ylthio)methyl thiocyanate; TCMTB	244-445-0	21564-17-0	Acute Tox. 2 * Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H330 H302 H319 H315 H317 H400 H410	GHS06 GHS09 Dgr	H330 H302 H319 H315 H317 H410			
613-120-00-9	bioresmethrin; (5-bezylfur-3-yl)methyl(1R)-*trans*-2,2-dimethyl-3-(2-methylpropenyl)cyclopropanecarboylate	249-014-0	28434-01-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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613-121-00-4	chlorsulfuron (ISO); 2-chloro-N-[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino]carbonyl]benzenesulphonamide	265-268-5	64902-72-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-122-00-X	diclobutrazole (ISO); (*R**, *R**)-(±)-β-[(2,4-dichlorophenyl)methyl]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol; (2*RS*, 3*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*—1,2,4-triazol-1yl)pentan-3-ol	—	75736-33-3	Eye Irrit. 2 Aquatic Chronic 2	H319 H411	GHS07 GHS09 Wng	H319 H411			
613-123-00-5	5,6-dihydro-3*H*-imidazo[2,1-*c*]-1,2,4-dithiazole-3-thione; etem	251-684-4	33813-20-6	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-124-00-0	fenpropimorph (ISO); *cis*-4-[3-(*p*-*tert*-butylphenyl)-2-methylpropyl]-2,6-dimethylmorpholine	266-719-9	67564-91-4	Repr. 2 Acute Tox. 4 * Skin Irrit. 2 Aquatic Chronic 2	H361d *** H302 H315 H411	GHS08 GHS07 GHS09 Wng	H361d *** H302 H315 H411			
613-125-00-6	hexythiazox(ISO); *trans*-5-(4-chlorophenyl)-*N*-cyclohexyl-4-methyl-2-oxo-3-thiazolidine-carboxamide	—	78587-05-0	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-126-00-1	imazapyr (ISO); 2-[4,5-dihydro-4-methyl-4-(1-methylethyl)-5-oxo-1*H*-imidazol-2-yl]-3-pyridine carboxylate	—	81334-34-1	Eye Irrit. 2 Aquatic Chronic 3	H319 H412	GHS07 Wng	H319 H412			
613-127-00-7	1,1-dimethylpiperidinium chloride; mepiquat chloride	246-147-6	24307-26-4	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			

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613-128-00-2	prochloraz (ISO); *N*-propyl-*N*-[2-(2,4,6-trichlorophenoxy) ethyl]-1*H*-imidazole-1-carboxamide	266-994-5	67747-09-5	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-129-00-8	metamitron (ISO); 4-amino-3-methyl-6-phenyl-1,2,4-triazin-5-one	255-349-3	41394-05-2	Acute Tox. 4 * Aquatic Acute 1	H302 H400	GHS07 GHS09 Wng	H302 H400			
613-131-00-9	pyroquilon (ISO); 1,2,5,6-tetrahydropyrrolo[3,2,1-ij]quinolin-4-one	—	57369-32-1	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
613-132-00-4	hexazinone (ISO); 3-cyclohexyl-6-dimethylamino-1-methyl-1,2,3,4-tetrahydro-1,3,5-triazine-2,4-dione	257-074-4	51235-04-2	Acute Tox. 4 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H302 H319 H400 H410	GHS07 GHS09 Wng	H302 H319 H410			
613-133-00-X	etridiazole (ISO); 5-ethoxy-3-trichloromethyl-1,2,4-thiadiazole	219-991-8	2593-15-9	Carc. 2 Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H351 H331 H312 H302 H400 H410	GHS06 GHS08 GHS09 Dgr	H351 H331 H312 H302 H410			
613-134-00-5	myclobutanil(ISO); 2-(4-chlorophenyl)-2-(1*H*-1,2,4-triazol-1-ylmethyl)hexanenitrile	—	88671-89-0	Repr. 2 Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 2	H361d *** H302 H319 H411	GHS08 GHS07 GHS09 Wng	H361d *** H302 H319 H411			
613-135-00-0	di(benzothiazol-2-yl) disulphide	204-424-9	120-78-5	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410	EUH031		

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613-136-00-6	*N*-cyclohexylbenzothiazole-2-sulphenamide	202-411-2	95-33-0	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
613-137-00-1	methabenzthiazuron (ISO); 1-(1,3-benzothiazol-2-yl)1,3-dimethylurea	242-505-0	18691-97-9	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-138-00-7	quinoxyfen (ISO); 5,7-dichloro-4-(4-fluorophenoxy)quinoline	—	124495-18-7	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
613-139-00-2	metsulfuron-methyl; 2-(4-methoxy-6-methyl-1,3,5-triazin-2-ylcarbamoylsulfamoyl) benzoic acid	—	74223-64-6	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-140-00-8	cycloheximide (ISO); 4-{(2*R*)-2-[(1*S*,3*S*,5*S*)-3,5-dimethyl-2-oxocyclohexyl]-2-hydroxyethyl}piperidine-2,6-dione	200-636-0	66-81-9	Muta. 2 Repr. 1B Acute Tox. 2 * Aquatic Chronic 2	H341 H360-D *** H300 H411	GHS06 GHS08 GHS09 Dgr	H341 H360D *** H300 H411			
613-141-00-3	1,4-diamino-2-(2-butyltetrazol-5-yl)-3-cyanoanthraquinone	401-470-3	93686-63-6	Aquatic Chronic 4	H413	—	H413			
613-142-00-9	*trans*—*N*-methyl-2-styryl-[4'-aminomethine-(1-acetyl-1-(2-methoxyphenyl)acetamido)] pyridinium acetate	405-860-4	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-143-00-4	1-(3-phenylpropyl)-2-methylpyridinium bromide	405-930-4	10551-42-5	Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 3	H302 H319 H412	GHS07 Wng	H302 H319 H412			
613-144-00-X	reaction products of: poly(vinyl acetate), partially hydrolyzed, with (*E*)-2-(4-formylstyryl)-3,4-dimethylthiazoliummethyl sulfate	406-460-2	125139-08-4	Aquatic Chronic 3	H412	—	H412			

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613-145-00-5	(*S*)-3-benzyloxycarbonyl-1,2,3,4-tetrahydroisoquinolinium 4-methylbenzenesulfonate	406-960-0	77497-97-3	Aquatic Chronic 2	H411	GHS09	H411			
613-146-00-0	*N*-ethyl-*N*-methylpiperidinium iodide	407-780-5	4186-71-4	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
613-147-00-6	4-[2-(1-methyl-2-(4-morpholinyl)ethoxy) ethyl]morpholine	407-940-4	111681-72-2	Eye Dam. 1	H318	GHS05 Dgr	H318			
613-148-00-1	tetrasodium 1,2-bis(4-fluoro-6-[5-(1-amino-2-sulfonatoanthrachinon-4-ylamino)-2,4,6-trimethyl-3-sulfonatophenylamino]-1,3,5-triazin-2-ylamino)ethane	411-240-4	143683-23-2	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
613-149-00-7	pyridaben (ISO); 2-*tert*-butyl-5-(4-*tert*-butylbenzylthio)-4-chloropyridazin-3(2*H*)-one	405-700-3	96489-71-3	Acute Tox. 3 * Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H331 H301 H400 H410	GHS06 GHS09 Dgr	H331 H301 H410			
613-150-00-2	2,2'-[3,3'-(piperazine-1,4-diyl)dipropyl]bis (1*H*-benzimidazo[2,1—*b*]benzo[*l*,*m*,*n*][3,8] phenanthroline-1,3,6-trione	406-295-6	—	Aquatic Chronic 4	H413	—	H413			
613-151-00-8	1-(3-mesyloxy-5-trityloxymethyl-2-D-threo-furyl)thymine	406-360-9	104218-44-2	Aquatic Chronic 4	H413	—	H413			
613-152-00-3	phenyl *N*-(4,6-dimethoxypyrimidin-2-yl)carbamate	406-600-2	89392-03-0	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-153-00-9	2,3,5-trichloropyridine	407-270-2	16063-70-0	Aquatic Chronic 3	H412	—	H412			
613-154-00-4	2-amino-4-chloro-6-methoxypyrimidine	410-050-9	5734-64-5	Acute Tox. 4 *	H302	GHS07 Wng	H302			

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613-155-00-X	5-chloro-2,3-difluoropyridine	410-090-7	89402-43-7	Flam. Liq. 3 Acute Tox. 4 * Aquatic Chronic 3	H226 H302 H412	GHS02 GHS07 Wng	H226 H302 H412			
613-156-00-5	2-butyl-4-chloro-5-formylimidazole	410-260-0	83857-96-9	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-157-00-0	2,4-diamino-5-methoxymethylpyrimidine	410-330-0	54236-98-5	Acute Tox. 4 * STOT RE 2 * Eye Irrit. 2	H302 H373 ** H319	GHS08 GHS07 Wng	H302 H373 ** H319			
613-158-00-6	2,3-dichloro-5-trifluoromethyl-pyridine	410-340-5	69045-84-7	Acute Tox. 4 * Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H332 H302 H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H332 H302 H318 H317 H411			
613-159-00-1	fenazaquin (ISO); 4-[2-[4-(1,1-dimethylethyl)phenyl]-ethoxy] quinazoline	410-580-0	120928-09-8	Acute Tox. 3 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H301 H332 H400 H410	GHS06 GHS09 Dgr	H301 H332 H410			
613-160-00-7	(1*S*)-2-methyl-2,5-diazobicyclo[2.2.1]heptane dihydrobromide	411-000-9	125224-62-6	Skin Sens. 1	H317	GHS07 Wng	H317			
613-163-00-3	azimsulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-[1-methyl-4-(2-methyl-2*H*-tetrazol-5-yl)pyrazol-5-ylsulfonyl]urea	—	120162-55-2	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-164-00-9	flufenacet (ISO); *N*-(4-fluorophenyl)-*N*-isopropyl-2-(5-trifluoromethyl-[1,3,4]thiadiazol-2-yloxy)acetamide	—	142459-58-3	Acute Tox. 4 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H373 ** H317 H400 H410	GHS08 GHS07 GHS09 Wng	H302 H373 ** H317 H410			
613-165-00-4	flupyrsulfuron-methyl-sodium (ISO); methyl 2-[[(4,6-dimethoxypyrimidin-2-ylcarbamoyl)sulfamoyl]-6-trifluoromethyl]nicotinate, monosodium salt	—	144740-54-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-166-00-X	flumioxazin (ISO); N-(7-fluoro-3,4-dihydro-3-oxo-4-prop-2-ynyl-2*H*—1,4-benzoxazin-6-yl)cyclohex-1-ene-1,2-dicarboxamide	—	103361-09-7	Repr. 1B Aquatic Acute 1 Aquatic Chronic 1	H360-D *** H400 H410	GHS08 GHS09 Dgr	H360D *** H410			
613-167-00-5	reaction mass of: 5-chloro-2-methyl-4-isothiazolin-3-one [EC no. 247-500-7] and 2-methyl-2*H* -isothiazol-3-one [EC no. 220-239-6] (3:1); reaction mass of: 5-chloro-2-methyl-4-isothiazolin-3-one [EC no. 247-500-7] and 2-methyl-4-isothiazolin-3-one [EC no. 220-239-6] (3:1)	—	55965-84-9	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H311 H301 H314 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H331 H311 H301 H314 H317 H410		Skin Corr. 1B; H314: C ≥ 0,6 % Skin Irrit. 2; H315: 0,06 % ≤ C < 0,6 % Eye Irrit. 2; H319: 0,06 % ≤ C < 0,6 % Skin Sens. 1; H317: C ≥ 0,0015 %	
613-168-00-0	1-vinyl-2-pyrrolidone	201-800-4	88-12-0	Carc. 2 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * STOT SE 3 Eye Dam. 1	H351 H332 H312 H302 H373 ** H335 H318	GHS06 GHS05 GHS09 Dgr	H351 H332 H312 H302 H373 ** H335 H318			D

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-169-00-6	9-vinylcarbazole	216-055-0	1484-13-5	Muta. 2 Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H341 H312 H302 H315 H317 H400 H410	GHS08 GHS07 GHS09 Wng	H341 H312 H302 H315 H317 H410			
613-170-00-1	2,2-ethylmethylthiazolidine	404-500-3	694-64-4	Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H302 H318 H317 H411			
613-171-00-7	hexaconazole (ISO); (*RS*)-2-(2,4-dichlorophenyl)-1-(1*H*-1,2,4-triazol-1-yl)hexan-2-ol	413-050-7	79983-71-4	Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H302 H317 H411	GHS07 GHS09 Wng	H302 H317 H411			
613-172-00-2	5-chloro-1,3-dihydro-2*H*-indol-2-one	412-200-9	17630-75-0	Repr. 2 Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 3	H361f *** H302 H317 H412	GHS08 GHS07 Wng	H361f *** H302 H317 H412			
613-173-00-8	fluquinconazole (ISO); 3-(2,4-dichlorophenyl)-6-fluoro-2-(1*H*-1,2,4-triazol-1-yl)quinazolin-4-(3*H*)-one	411-960-9	136426-54-5	Acute Tox. 3 * Acute Tox. 3 * STOT RE 1 Acute Tox. 4 * Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H331 H301 H372 ** H312 H315 H400 H410	GHS06 GHS08 GHS09 Dgr	H331 H301 H372 ** H312 H315 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-174-00-3	(±) 2-(2,4-dichlorophenyl)-3-(1*H*-1,2,4-triazol-1-yl)propyl-1,1,2,2-tetrafluoroethylether	407-760-6	112281-77-3	Carc. 2 Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 2	H351 H332 H302 H411	GHS08 GHS07 GHS09 Wng	H351 H332 H302 H411			
613-175-00-9	epoxiconazole(ISO); (2*RS*,3*SR*)-3-(2-chlorophenyl)-2-(4-fluorophenyl)-[(1*H*-1,2,4-triazol-1-yl)methyl]oxirane	406-850-2	133855-98-8	Carc. 2 Repr. 2 Aquatic Chronic 2	H351 H361f-d H411	GHS08 GHS09 Wng	H351 H361fd H411			
613-176-00-4	2-methyl-2-azabicyclo[2.2.1]heptane	404-810-9	4524-95-2	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B	H226 H312 H302 H373 ** H314	GHS02 GHS08 GHS05 GHS07 Dgr	H226 H312 H302 H373 ** H314			
613-177-00-X	8-amino-7-methylquinoline	412-760-4	5470-82-6	Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Chronic 2	H312 H302 H317 H411	GHS07 GHS09 Wng	H312 H302 H317 H411			
613-178-00-5	4-ethyl-2-methyl-2-isopentyl-1,3-oxazolidine	410-470-2	137796-06-6	Skin Corr. 1B Skin Sens. 1	H314 H317	GHS05 GHS07 Dgr	H314 H317		STOT SE 3; H335: C ≥ 5 %	
613-179-00-0	lithium 3-oxo-1,2(2*H*)-benzisothiazol-2-ide	411-690-1	111337-53-2	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Chronic 2	H302 H314 H317 H411	GHS05 GHS07 Dgr	H302 H314 H317 H411			
613-180-00-6	*N*-(1,1-dimethylethyl)bis(2-benzothiazolesulfen)amide	407-430-1	3741-80-8	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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613-181-00-1	5,5-dimethyl-perhydro-pyrimidin-2-one α-(4-trifluoromethylstyryl)-α-(4-trifluoromethyl)cinnamylidenehydrazone	405-090-9	67485-29-4	STOT RE 1 Acute Tox. 4 * Eye Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H372 ** H302 H319 H400 H410	GHS08 GHS07 GHS09 Dgr	H372 ** H302 H319 H410			
613-182-00-7	1-(1-naphthylmethyl)quinolinium chloride	406-220-7	65322-65-8	Carc. 2 Muta. 2 Acute Tox. 4 * Skin Irrit. 2 Eye Dam. 1 Aquatic Chronic 3	H351 H341 H302 H315 H318 H412	GHS08 GHS05 GHS07 Dgr	H351 H341 H302 H315 H318 H412			
613-183-00-2	reaction mass of: 5-(*N*-methylperfluorooctylsulfonamido)methyl-3-octadecyl-1,3-oxazolidin-2-one; 5-(*N*-methylperfluoroheptylsulfonamido) methyl-3-octadecyl-1,3-oxazolidin-2-one	413-640-4	—	STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H373 ** H400 H410	GHS08 GHS09 Wng	H373 ** H410			
613-184-00-8	nitrilotriethyleneammoniopropane-2-ol 2-ethylhexanoate	413-670-8	—	Eye Irrit. 2 Skin Sens. 1	H319 H317	GHS07 Wng	H319 H317			
613-185-00-3	2,3,5,6-tetrahydro-2-methyl-2*H*-cyclopenta[*d*]-1,2-thiazol-3-one	407-630-9	82633-79-2	Acute Tox. 3 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H301 H318 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H301 H318 H317 H410			
613-186-00-9	(2*R*,3*R*)-3-((*R*)-1-(*tert*-butyldimethylsiloxy) ethyl)-4-oxoazetidin-2-yl acetate	408-050-9	76855-69-1	Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H319 H317 H411	GHS07 GHS09 Wng	H319 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-188-00-X	1-(3-(4-fluorophenoxy)propyl)-3-methoxy-4-piperidinone	411-500-7	116256-11-2	Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H302 H318 H317 H411			
613-189-00-5	1,4,7,10-tetrakis(*p*-toluensulfonyl)-1,4,7,10-tetraazacyclododecane	414-030-0	52667-88-6	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
613-190-00-0	disodium 1-amino-4-(2-(5-chloro-6-fluoro-pyrimidin-4-ylamino-methyl)-4-methyl-6-sulfo-phenylamino)-9,10-dioxo-9,10-dihydro-anthracene-2-sulfonate	414-040-5	149530-93-8	Acute Tox. 4 * Skin Sens. 1	H302 H317	GHS07 Wng	H302 H317			
613-191-00-6	3-ethyl-2-methyl-2-(3-methylbutyl)-1,3-oxazolidine	421-150-7	143860-04-2	Repr. 1B Skin Corr. 1B Aquatic Acute 1 Aquatic Chronic 1	H360F *** H314 H400 H410	GHS08 GHS05 GHS09 Dgr	H360F *** H314 H410			
613-193-00-7	pentakis[3-(dimethylammonio)propylsulfamoyl]-[(6-hydroxy-4,4,8,8-tetramethyl-4,8-diazoniaundecane-1,11-diyldisulfamoyl)di[phthalocyaninecopper(II)]] heptalactate	414-930-3	—	Aquatic Chronic 2	H411	GHS09	H411			
613-194-00-2	6,13-dichloro-3,10-bis{2-[4-fluoro-6-(2-sulfophenylamino)-1,3,5-triazin-2-ylamino]propylamino}benzo[5,6][1,4]oxazino[2,3-.b.]phenoxazine-4,11-disulphonic acid, lithium-, sodium salt	418-000-8	163062-28-0	Eye Dam. 1	H318	GHS05 Dgr	H318			
613-195-00-8	2,2-(1,4-phenylene)bis((4*H*—3,1-benzoxazine-4-one)	418-280-1	18600-59-4	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-196-00-3	5-[[4-chloro-6-[[2-[[4-fluoro-6-[[5-hydroxy-6-[(4-methoxy-2-sulfophenyl)azo]-7-sulfo-2-naphthalenyl]amino]-1,3,5-triazin-2-yl]amino]-1-methylethyl]amino]-1,3,5-triazin-2-yl]amino]-3-[[4-(ethenylsulfonyl)phenyl]azo]-4-hydroxy-naphtalene-2,7-disulfonic acid, sodium salt	418-380-5	168113-78-8	Eye Dam. 1	H318	GHS05 Dgr	H318			
613-197-00-9	reaction mass of: 2,4,6-tri(butylcarbamoyl)-1,3,5-triazine; 2,4,6-tri(methylcarbamoyl)-1,3,5-triazine; [(2-butyl-4,6-dimethyl)tricarbamoyl]-1,3,5-triazine; [(2,4-dibutyl-6-methyl)tricarbamoyl]-1,3,5-triazine	420-390-1	187547-46-2	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-199-00-X	reaction mass of: 1,3,5-tris(3-aminomethylphenyl)-1,3,5-(1*H*,3*H*,5*H*)-triazine-2,4,6-trione; reaction mass of oligomers of 3,5-bis(3-aminomethylphenyl)-1-poly[3,5-bis(3-aminomethylphenyl)-2,4,6-trioxo-1,3,5-(1*H*,3*H*,5*H*)-triazin-1-yl]-1,3,5-(1*H*,3*H*,5*H*)-triazine-2,4,6-trione	421-550-1	—	Carc. 1B Repr. 1B Skin Sens. 1 Aquatic Chronic 3	H350 H360-D *** H317 H412	GHS08 Dgr	H350 H360D *** H317 H412			
613-200-00-3	Reaction product of: copper, (29*H*,31*H*-phthalocyaninato(2-)-*N*29,*N*30,*N*31,*N*32)-, chlorosulfuric acid and 3-(2-sulfooxyethylsulfonyl)aniline, sodium salts	420-980-7	—	Eye Dam. 1	H318	GHS05 Dgr	H318			
613-201-00-9	(*R*)-5-bromo-3-(1-methyl-2-pyrrolidinyl methyl)-1*H*-indole	422-390-5	143322-57-0	Repr. 2 STOT RE 1 Acute Tox. 4 * Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361f *** H372 ** H332 H302 H317 H400 H410	GHS08 GHS07 GHS09 Dgr	H361f *** H372 ** H332 H302 H317 H410	EUH070		

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-202-00-4	pymetrozine (ISO); (*E*)-4,5-dihydro-6-methyl-4-(3-pyridylmethyleneamino)-1,2,4-triazin-3(2*H*)-one	—	123312-89-0	Carc. 2 Aquatic Chronic 3	H351 H412	GHS08 Wng	H351 H412			
613-203-00-X	pyraflufen-ethyl; [1] pyraflufen [2]	-[1] -[2]	129630-19-9 [1] 129630-17-7 [2]	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-204-00-5	oxadiargyl (ISO); 3-[2,4-dichloro-5-(2-propynyloxy)phenyl]-5-(1,1-dimethylethyl)-1,3,4-oxadiazol-2(3*H*)-one; 5-*tert*-butyl-3-[2,4-dichloro-5-(prop-2-ynyloxy)phenyl]-1,3,4-oxadiazol-2(3*H*)-one	254-637-6	39807-15-3	Repr. 2 STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H361d *** H373 ** H400 H410	GHS08 GHS09 Wng	H361d *** H373 ** H410			
613-205-00-0	propiconazole(ISO); (±) 1-[2-(2,4-dichlorophenyl)-4-propyl-1,3-dioxolan-2-ylmethyl]-1*H*—1,2,4-triazole	262-104-4	60207-90-1	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
613-206-00-6	fenamidone (ISO); (*S*)-5-methyl-2-methylthio-5-phenyl-3-phenylamino-3,5-dihydroimidazol-4-one	—	161326-34-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-208-00-7	imazamox (ISO); (*RS*)-2-(4-isopropyl-4-methyl-5-oxo-2-imidazolin-2-yl)-5-methoxymethylnicotinic acid	—	114311-32-9	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-209-00-2	*cis*-1-(3-chloropropyl)-2,6-dimethyl-piperidin hydrochloride	417-430-3	63645-17-0	Acute Tox. 3 * STOT RE 2 * Skin Sens. 1 Aquatic Chronic 2	H301 H373 ** H317 H411	GHS06 GHS08 GHS09 Dgr	H301 H373 ** H317 H411			
613-210-00-8	2-(3-chloropropyl)-2,5,5-trimethyl-1,3-dioxane	417-650-1	88128-57-8	STOT RE 2 * Aquatic Chronic 3	H373 ** H412	GHS08 Wng	H373 ** H412			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
613-211-00-3	*N*-methyl-4-(*p*-formylstyryl)pyridinium methylsulfate	418-240-3	74401-04-0	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
613-212-00-9	4-[4-(2-ethylhexyloxy)phenyl](1,4-thiazinane-1,1-dioxide)	418-320-8	133467-41-1	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
613-213-00-4	*cis*-1-benzoyl-4-[(4-methylsulfonyl)oxy]-L-proline	416-040-0	120807-02-5	Aquatic Chronic 3	H412	—	H412			
613-214-00-X	*N*,*N*-di-*n*-butyl-2-(1,2-dihydro-3-hydroxy-6-isopropyl-2-quinolylidene)-1,3-dioxoindan-5-carboxamide	416-260-7	147613-95-4	Aquatic Chronic 4	H413	—	H413			
613-215-00-5	2-chloromethyl-3,4-dimethoxypyridinium chloride	416-440-5	72830-09-2	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H312 H302 H373 ** H315 H318 H317 H411	GHS08 GHS05 GHS07 GHS09 Dgr	H312 H302 H373 ** H315 H318 H317 H411			
613-216-00-0	6-*tert*-butyl-7-(6-diethylamino-2-methyl-3-pyridylimino)-3-(3-methylphenyl)pyrazolo[3,2-c][1,2,4]triazole	416-490-8	162208-01-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-217-00-6	4-[3-(3,5-di-*tert*-butyl-4-hydroxyphenyl)propionyloxy]-1-[2-[3-(3,5-di-*tert*-butyl-4-hydrophenyl)propionyloxy]ethyl]-2,2,6,6-tetramethylpiperidine	416-770-1	73754-27-5	Aquatic Chronic 4	H413	—	H413			
613-218-00-1	6-hydroxyindole	417-020-4	2380-86-1	Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H302 H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H302 H318 H317 H411			

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613-219-00-7	7a-ethyl-3,5-bis(1-methylethyl)-2,3,4,5-tetrahydrooxazolo[3,4-c]-2,3,4,5-tetrahydrooxazole	417-140-7	79185-77-6	Skin Irrit. 2 Aquatic Chronic 2	H315 H411	GHS07 GHS09 Wng	H315 H411			
613-220-00-2	trans-(4*S*,6*S*)-5,6-dihydro-6-methyl-4*H*-thieno[2,3-b]thiopyran-4-ol, 7,7-dioxide	417-290-3	147086-81-5	Acute Tox. 4 *	H302	GHS07 Wng	H302			
613-221-00-8	2-chloro-5-methyl-pyridine	418-050-0	18368-64-4	Acute Tox. 4 * Acute Tox. 4 * Skin Irrit. 2 Aquatic Chronic 3	H312 H302 H315 H412	GHS07 Wng	H312 H302 H315 H412			
613-222-00-3	4-(1-oxo-2-propenyl)-morpholine	418-140-1	5117-12-4	Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1	H302 H373 ** H318 H317	GHS08 GHS05 GHS07 Dgr	H302 H373 ** H318 H317			
613-223-00-9	*N*-isopropyl-3-(4-fluorophenyl)-1*H*-indole	418-790-4	93957-49-4	Aquatic Chronic 4	H413	—	H413			
613-224-00-4	2,5-dimercaptomethyl-1,4-dithiane	419-770-8	136122-15-1	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H314 H317 H410			
613-225-00-X	reaction mass of:[2-(anthraquinon-1-ylamino)-6-[(5-benzoylamino)-anthraquinone-1-ylamino]-4-phenyl]-1,3,5-triazine; 2,6-bis-[(5-benzoylamino)-anthraquinon-1-ylamino]-4-phenyl-1,3,5-triazine.	421-290-9	—	STOT RE 2 * Aquatic Chronic 4	H373 ** H413	GHS08 Wng	H373 ** H413			
613-226-00-5	1-(2-(ethyl(4-(4-(4-(4-(ethyl(2-pyridinoethyl)amino)-2-methylphenylazo)benzoylamino)-phenylazo)-3-methylphenyl)amino)ethyl)-pyridinium dichloride	420-950-3	163831-67-2	Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H318 H400 H410	GHS05 GHS09 Dgr	H318 H410			

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613-227-00-0	(±)-[(*R**,*R**) and (*R**,*S**)]-6-fluoro-3,4-dihydro-2-oxiranyl-2*H*-1-benzopyran	419-600-2	99199-90-3	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
613-228-00-6	(±)-(*R**,*S**)-6-fluoro-3,4-dihydro-2-oxiranyl-2*H*-1-benzopyran	419-630-6	793669-26-8	Aquatic Chronic 2	H411	GHS09	H411			
613-230-00-7	florasulam (ISO); 2',6',8-trifluoro-5-methoxy-5-triazolo [1,5-c]; pyrimidine-2-sulfonanilide	—	145701-23-1	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
613-233-00-3	4,4'-(oxy-(bismethylene))-bis-1,3-dioxolane	423-230-7	56552-15-9	Eye Dam. 1	H318	GHS05 Dgr	H318			
614-001-00-4	nicotine (ISO); 3-(N-methyl-2-pyrrolidinyl)pyridine	200-193-3	54-11-5	Acute Tox. 1 Acute Tox. 3 * Aquatic Chronic 2	H310 H301 H411	GHS06 GHS09 Dgr	H310 H301 H411			
614-002-00-X	salts of nicotine	—	—	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 * Aquatic Chronic 2	H330 H310 H300 H411	GHS06 GHS09 Dgr	H330 H310 H300 H411			A
614-003-00-5	strychnine	200-319-7	57-24-9	Acute Tox. 1 Acute Tox. 2 * Aquatic Acute 1 Aquatic Chronic 1	H310 H300 H400 H410	GHS06 GHS09 Dgr	H310 H300 H410			
614-004-00-0	salts of strychnine	—	—	Acute Tox. 2 * Acute Tox. 2 * Aquatic Acute 1 Aquatic Chronic 1	H330 H300 H400 H410	GHS06 GHS09 Dgr	H330 H300 H410			A

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
614-005-00-6	colchicine	200-598-5	64-86-8	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-006-00-1	brucine; 2,3-dimethoxystrychnine	206-614-7	357-57-3	Acute Tox. 2 * Acute Tox. 2 * Aquatic Chronic 3	H330 H300 H412	GHS06 Dgr	H330 H300 H412			
614-007-00-7	brucine sulphate; [1] brucine nitrate; [2] strychnidin-10-one, 2,3-dimethoxy-, mono [(R)-1-methylheptyl 1,2-benzenedicarboxylate]; [3] strychnidin-10-one, 2,3-dimethoxy-, compd. with (S)mono(1-methylheptyl)-1,2-benzenedicarboxylate (1:1) [4]	225-432-9 [1] 227-317-9 [2] 269-439-5 [3] 269-710-8 [4]	4845-99-2 [1] 5786-97-0 [2] 68239-26-9 [3] 68310-42-9 [4]	Acute Tox. 2 * Acute Tox. 2 * Aquatic Chronic 3	H330 H300 H412	GHS06 Dgr	H330 H300 H412			A
614-008-00-2	aconitine	206-121-7	302-27-2	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-009-00-8	salts of aconitine	—	—	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			A
614-010-00-3	atropine	200-104-8	51-55-8	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-011-00-9	salts of atropine	—	—	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			A
614-012-00-4	hyoscyamine	202-933-0	101-31-5	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-013-00-X	salts of hyoscyamine	—	—	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			A
614-014-00-5	hyoscine	200-090-3	51-34-3	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 *	H330 H310 H300	GHS06 Dgr	H330 H310 H300			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
614-015-00-0	salts of hyoscine	—	—	Acute Tox. 2 * Acute Tox. 1 Acute Tox. 2 *	H330 H310 H300	GHS06 Dgr	H330 H310 H300			A
614-016-00-6	pilocarpine	202-128-4	92-13-7	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-017-00-1	salts of pilocarpine	—	—	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			A
614-018-00-7	papaverine	200-397-2	58-74-2	Acute Tox. 4 *	H302	GHS07 Wng	H302			
614-019-00-2	salts of papaverine	—	—	Acute Tox. 4 *	H302	GHS07 Wng	H302			A
614-020-00-8	physostigmine	200-332-8	57-47-6	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			
614-021-00-3	salts of physostigmine	—	—	Acute Tox. 2 * Acute Tox. 2 *	H330 H300	GHS06 Dgr	H330 H300			A
614-022-00-9	digitoxin	200-760-5	71-63-6	Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H331 H301 H373 **	GHS06 GHS08 Dgr	H331 H301 H373 **			
614-023-00-4	ephedrine	206-080-5	299-42-3	Acute Tox. 4 *	H302	GHS07 Wng	H302			
614-024-00-X	salts of ephedrine	—	—	Acute Tox. 4 *	H302	GHS07 Wng	H302			A
614-025-00-5	ouabain	211-139-3	630-60-4	Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H331 H301 H373 **	GHS06 GHS08 Dgr	H331 H301 H373 **			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
614-026-00-0	strophantin-K	234-239-9	11005-63-3	Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 *	H331 H301 H373 **	GHS06 GHS08 Dgr	H331 H301 H373 **			
614-027-00-6	bufa-4,20,22-trienolide, 6-(acetyloxy)-3-(β-D-glucopyranosyloxy)-8,14-dihydroxy-, (3β, 6β)-; red squill; scilliroside	208-077-4	507-60-8	Acute Tox. 2 *	H300	GHS06 Dgr	H300			
614-028-00-1	reaction mass of: 2-ethylhexyl mono-D-glucopyranoside; 2-ethylhexyl di-D-glucopyranoside	414-420-0	—	Eye Dam. 1	H318	GHS05 Dgr	H318			
614-029-00-7	constitutional isomers of penta-O-allyl-β-D-fructofuranosyl-α-D-glucopyranoside; constitutional isomers of hexa-O-allyl-β-D-fructofuranosyl-α-D-glucopyranoside; constitutional isomers of hepta-O-allyl-β-D-fructofuransoyl-α-D-glucopyranoside	419-640-0	68784-14-5	Acute Tox. 4 *	H302	GHS07 Wng	H302			
615-001-00-7	methyl isocyanate	210-866-3	624-83-9	Flam. Liq. 2 Repr. 2 Acute Tox. 2 * Acute Tox. 3 * Acute Tox. 3 * STOT SE 3 Skin Irrit. 2 Eye Dam. 1 Resp. Sens. 1 Skin Sens. 1	H225 H361d *** H330 H311 H301 H335 H315 H318 H334 H317	GHS02 GHS06 GHS06 GHS08 GHS05 Dgr	H225 H361d *** H330 H311 H301 H335 H315 H318 H334 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
615-002-00-2	methyl isothiocyanate	209-132-5	556-61-6	Acute Tox. 3 * Acute Tox. 3 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H301 H314 H317 H400 H410	GHS06 GHS05 GHS09 Dgr	H331 H301 H314 H317 H410			
615-003-00-8	thiocyanic acid	207-337-4	463-56-9	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H332 H312 H302 H412	GHS07 Wng	H332 H312 H302 H412	EUH032		
615-004-00-3	salts of thiocyanic acid	—	—	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H332 H312 H302 H412	GHS07 Wng	H332 H312 H302 H412	EUH032		A
615-005-00-9	4,4'-methylenediphenyl diisocyanate; diphenylmethane-4,4'-diisocyanate; [1] 2,2'-methylenediphenyl diisocyanate; diphenylmethane-2,2'-diisocyanate; [2] *o*-(*p*-isocyanatobenzyl)phenyl isocyanate; diphenylmethane-2,4'-diisocyanate; [3] methylenediphenyl diisocyanate [4]	202-966-0 [1] 219-799-4 [2] 227-534-9 [3] 247-714-0 [4]	101-68-8 [1] 2536-05-2 [2] 5873-54-1 [3] 26447-40-5 [4]	Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Skin Sens. 1	H332 H319 H335 H315 H334 H317	GHS08 GHS07 Dgr	H332 H319 H335 H315 H334 H317		Eye Irrit.; H319: C ≥ 5 % STOT SE 3; H335: C ≥ 5 % Skin Irrit. 2; H315: C ≥ 5 % Resp. Sens. 1; H334: C ≥ 0,1 %	C 2
615-006-00-4	2-methyl-*m*-phenylene diisocyanate; toluene-2,4-di-isocyanate; [1] 4-methyl-*m*-phenylene diisocyanate; toluene-2,6-di-isocyanate; [2] *m*-tolylidene diisocyanate; toluene-diisocyanate [3]	202-039-0 [1] 209-544-5 [2] 247-722-4 [3]	91-08-7 [1] 584-84-9 [2] 26471-62-5 [3]	Carc. 2 Acute Tox. 2 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 3	H351 H330 H319 H335 H315 H334 H317 H412	GHS06 GHS08 Dgr	H351 H330 H319 H335 H315 H334 H317 H412		Resp. Sens. 1; H334: C ≥ 0,1 %	C

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
615-007-00-X	1,5-naphthylene diisocyanate	221-641-4	3173-72-6	Acute Tox. 4 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Aquatic Chronic 3	H332 H319 H335 H315 H334 H412	GHS08 GHS07 Dgr	H332 H319 H335 H315 H334 H412			
615-008-00-5	3-isocyanatomethyl-3,5,5-trimethylcyclohexyl isocyanate; isophorone di-isocyanate	223-861-6	4098-71-9	Acute Tox. 3 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 2	H331 H319 H335 H315 H334 H317 H411	GHS06 GHS08 GHS09 Dgr	H331 H319 H335 H315 H334 H317 H411		* Resp. Sens. 1; H334: C ≥ 0,5 % Skin Sens.1; H317: C ≥ 0,5 %	2
615-009-00-0	4,4'-methylenedi(cyclohexyl isocyanate); dicyclohexylmethane-4,4'-di-isocyanate	225-863-2	5124-30-1	Acute Tox. 3 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Skin Sens. 1	H331 H319 H335 H315 H334 H317	GHS06 GHS08 Dgr	H331 H319 H335 H315 H334 H317		* Resp. Sens. 1; H334: C ≥ 0,5 % Skin Sens. 1; H317: C ≥ 0,5 %	2
615-010-00-6	2,2,4-trimethylhexamethylene-1,6-di-isocyanate; [1] 2,4,4-trimethylhexamethylene-1,6-di-isocyanate [2]	241-001-8 [1] 239-714-4 [2]	16938-22-0 [1] 15646-96-5 [2]	Acute Tox. 3 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H331 H319 H335 H315 H334	GHS06 GHS08 Dgr	H331 H319 H335 H315 H334		* Resp. Sens. 1; H334: C ≥ 0,5 % Skin Sens. 1; H317: C ≥ 0,5 %	C 2
615-011-00-1	hexamethylene-di-isocyanate	212-485-8	822-06-0	Acute Tox. 3 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1 Skin Sens. 1	H331 H319 H335 H315 H334 H317	GHS06 GHS08 Dgr	H331 H319 H335 H315 H334 H317		* Resp. Sens. 1; H334: C ≥ 0,5 % Skin Sens. 1; H317: C ≥ 0,5 %	2

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
615-012-00-7	4-isocyanatosulphonyltoluene; tosyl isocyanate	223-810-8	4083-64-1	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334	EUH014	Eye Irrit.; H319: C ≥ 5 % STOT SE 3; H335: C ≥ 5 % Skin Irrit. 2; H315: C ≥ 5 %	
615-013-00-2	cyanamide; carbanonitril	206-992-3	420-04-2	Acute Tox. 3 * Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1	H301 H312 H319 H315 H317	GHS06 Dgr	H301 H312 H319 H315 H317			
615-014-00-8	tris(1-dodecyl-3-methyl-2-phenylbenzimidazolium)hexacyanoferrate	—	7276-58-6	Acute Tox. 4 *	H302	GHS07 Wng	H302			
615-015-00-3	1,7,7-trimethylbicyclo(2,2,1)hept-2-yl thiocyanatoacetate; isobornyl thiocyanoacetate	204-081-5	115-31-1	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
615-016-00-9	potassium cyanate	209-676-3	590-28-3	Acute Tox. 4 *	H302	GHS07 Wng	H302			
615-017-00-4	calcium cyanamide	205-861-8	156-62-7	Acute Tox. 4 * STOT SE 3 Eye Dam. 1	H302 H335 H318	GHS05 GHS07 Dgr	H302 H335 H318			
615-018-00-X	2-(2-butoxyethoxy)ethyl thiocyanate	203-985-7	112-56-1	Flam. Liq. 3 Acute Tox. 3 * Acute Tox. 3 *	H226 H311 H301	GHS02 GHS06 Dgr	H226 H311 H301			
615-019-00-5	dicyclohexylcarbodiimide	208-704-1	538-75-0	Acute Tox. 3 * Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1	H311 H302 H318 H317	GHS06 GHS05 Dgr	H311 H302 H318 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
615-020-00-0	methylene dithiocyanate	228-652-3	6317-18-6	Acute Tox. 2 * Acute Tox. 3 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1	H330 H301 H314 H317 H400	GHS06 GHS05 GHS09 Dgr	H330 H301 H314 H317 H400			
615-021-00-6	1,3,5-tris(oxiranylmethyl)-1,3,5-triazine-2,4,6(1*H*,3*H*,5*H*)-trione; TGIC	219-514-3	2451-62-9	Muta. 1B Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H340 H331 H301 H373 ** H318 H317 H412	GHS06 GHS08 GHS05 Dgr	H340 H331 H301 H373 ** H318 H317 H412			
615-022-00-1	methyl 3-isocyanatosulfonyl-2-thiophene-carboxylate	410-550-7	79277-18-2	STOT RE 2 * Resp. Sens. 1 Skin Sens. 1	H373 ** H334 H317	GHS08 Dgr	~~E; R2~~ H373 ** H334 H317	EUH014		
615-023-00-7	2-(isocyanatosulfonylmethyl)benzoic acid methyl ester; (alt.):methyl 2-(isocyanatosulfonylmethyl) benzoate	410-900-9	83056-32-0	Flam. Liq. 3 Muta. 2 Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Resp. Sens. 1	H226 H341 H332 H373 ** H318 H334	GHS02 GHS08 GHS05 GHS07 Dgr	H226 H341 H332 H373 ** H318 H334	EUH014		
615-024-00-2	2-phenylethylisocyanate	413-080-0	1943-82-4	Acute Tox. 3 * Acute Tox. 4 * Skin Corr. 1A Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 2	H331 H302 H314 H334 H317 H411	GHS06 GHS08 GHS05 GHS09 Dgr	H331 H302 H314 H334 H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
615-025-00-8	4,4'-ethylidenediphenyl dicyanate	405-740-1	47073-92-7	Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H302 H373 ** H318 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H332 H302 H373 ** H318 H410			
615-026-00-3	4,4'-methylenebis(2,6-dimethylphenyl cyanate)	405-790-4	101657-77-6	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
615-028-00-4	ethyl 2-(isocyanatosulfonyl)benzoate	410-220-2	77375-79-2	**** Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Resp. Sens. 1 Skin Sens. 1	H302 H373 ** H318 H334 H317	GHS08 GHS05 GHS07 Dgr	~~E; R2~~ H302 H373 ** H318 H334 H317	EUH014		
615-029-00-X	2,5-bis-isocyanatomethyl-bicyclo[2.2.1]heptane	411-280-2	—	Acute Tox. 2 * Acute Tox. 4 * Skin Corr. 1B Resp. Sens. 1 Skin Sens. 1 Aquatic Chronic 3	H330 H302 H314 H334 H317 H412	GHS06 GHS08 GHS05 Dgr	H330 H302 H314 H334 H317 H412			
615-030-00-5	alkali salts, alkali earth salts and other salts of thiocyanic acid not mentioned elsewhere in this Annex	—	—	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 3	H332 H312 H302 H412	GHS07 Wng	H332 H312 H302 H412	EUH032		A
615-031-00-0	thallium salt of thiocyanic acid	222-571-7	3535-84-0	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Chronic 2	H332 H312 H302 H411	GHS07 GHS09 Wng	H332 H312 H302 H411	EUH032		A

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615-032-00-6	metal salts of thiocyanic acid not mentioned elsewhere in this Annex	—	—	Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H332 H312 H302 H400 H410	GHS07 GHS09 Wng	H332 H312 H302 H410	EUH032		A
616-001-00-X	*N,N*-dimethylformamide; dimethyl formamide	200-679-5	68-12-2	Repr. 1B Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2	H360-D *** H332 H312 H319	GHS08 GHS07 Dgr	H360D *** H332 H312 H319			
616-002-00-5	2-fluoroacetamide	211-363-1	640-19-7	Acute Tox. 2 * Acute Tox. 3 *	H300 H311	GHS06 Dgr	H300 H311			
616-003-00-0	acrylamide; prop-2-enamide	201-173-7	79-06-1	Carc. 1B Muta. 1B Repr. 2 Acute Tox. 3 * STOT RE 1 Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1	H350 H340 H361f *** H301 H372 ** H332 H312 H319 H315 H317	GHS06 GHS08 Dgr	H350 H340 H361f *** H301 H372 ** H332 H312 H319 H315 H317			D
616-004-00-6	allidochlor (ISO); *N,N*-diallylchloroacetamide	202-270-7	93-71-0	Acute Tox. 4 * Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 2	H312 H302 H319 H315 H411	GHS07 GHS09 Wng	H312 H302 H319 H315 H411			
616-005-00-1	chlorthiamid (ISO); 2,6-dichloro (thiobenzamide)	217-637-7	1918-13-4	Acute Tox. 4 *	H302	GHS07 Wng	H302			

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616-006-00-7	dichlofluanid (ISO); *N*-dichlorofluoromethylthio-*N*',*N*'-dimethyl-*N*-phenylsulphamide	214-118-7	1085-98-9	Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H319 H317 H400 H410	GHS07 GHS09 Wng	H332 H319 H317 H410			
616-007-00-2	diphenamid (ISO); *N*,*N*-dimethyl-2,2-diphenylacetamide	213-482-4	957-51-7	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
616-008-00-8	propachlor (ISO); 2-chloro-*N*-isopropylacetanilide; α-chloro-*N*-isopropylacetanilide	217-638-2	1918-16-7	Acute Tox. 4 * Eye Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H319 H317 H400 H410	GHS07 GHS09 Wng	H302 H319 H317 H410			
616-009-00-3	propanil (ISO); 3',4'-dichloropropionanilide	211-914-6	709-98-8	Acute Tox. 4 * Aquatic Acute 1	H302 H400	GHS07 GHS09 Wng	H302 H400			
616-010-00-9	tosylchloramide sodium	204-854-7	127-65-1	Acute Tox. 4 * Skin Corr. 1B Resp. Sens. 1	H302 H314 H334	GHS08 GHS05 GHS07 Dg	H302 H314 H334	EUH031		
616-011-00-4	*N*,*N*-dimethylacetamide	204-826-4	127-19-5	Repr. 1B Acute Tox. 4 * Acute Tox. 4 *	H360-D *** H332 H312	GHS08 GHS07 Dgr	H360D *** H332 H312		Repr. 1B; H360D: C ≥ 5 %	
616-012-00-X	*N*-(dichlorofluoromethylthio)phthalimide; *N*-(fluorodichloromethylthio)phthalimide	211-952-3	719-96-0	Skin Irrit. 2	H315	GHS07 Wng	H315			
616-013-00-5	butyraldehyde oxime	203-792-8	110-69-0	Acute Tox. 3 * Acute Tox. 4 * Eye Irrit. 2	H311 H302 H319	GHS06 Dgr	H311 H302 H319			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-014-00-0	2-butanone oxime; ethyl methyl ketoxime; ethyl methyl ketone oxime	202-496-6	96-29-7	Carc. 2 Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1	H351 H312 H318 H317	GHS08 GHS05 GHS07 Dgr	H351 H312 H318 H317			
616-015-00-6	alachlor (ISO); 2-chloro-2',6'-diethyl-*N*-(methoxymethyl) acetanilide	240-110-8	15972-60-8	Carc. 2 Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H351 H302 H317 H400 H410	GHS08 GHS07 GHS09 Wng	H351 H302 H317 H410		M=10	
616-016-00-1	1-(3,4-dichlorophenylimino) thiosemicarbazide	—	5836-73-7	Acute Tox. 2 *	H300	GHS06 Dgr	H300			
616-017-00-7	cartap hydrochloride	239-309-2	15263-52-2	Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H312 H302 H400 H410	GHS07 GHS09 Wng	H312 H302 H410			
616-018-00-2	*N*,*N*-diethyl-*m*-toluamide; deet	205-149-7	134-62-3	Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 3	H302 H319 H315 H412	GHS07 Wng	H302 H319 H315 H412			
616-019-00-8	perfluidone (ISO); 1,1,1-trifluoro-*N*-(4-phenylsulphonyl-*o*-tolyl)methanesulphonamide;	253-718-3	37924-13-3	Acute Tox. 4 * Eye Irrit. 2	H302 H319	GHS07 Wng	H302 H319			
616-020-00-3	tebuthiuron (ISO); 1-(5-*tert*-butyl-1,3,4-thiadiazol-2-yl)-1,3-dimethylurea	251-793-7	34014-18-1	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-021-00-9	thiazafluron (ISO); 1,3-dimethyl-1-(5-trifluoromethyl-1,3,4-thiadiazol-2-yl)urea	246-901-4	25366-23-8	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
616-022-00-4	acetamide	200-473-5	60-35-5	Carc. 2	H351	GHS08 Wng	H351			
616-023-00-X	*N*-hexadecyl(or octadecyl)-*N*-hexadecyl(or octadecyl)benzamide	401-980-6	—	Skin Irrit. 2 Skin Sens. 1	H315 H317	GHS07 Wng	H315 H317			
616-024-00-5	2-(4,4-dimethyl-2,5-dioxooxazolidin-1-yl)-2-chloro-5-(2-(2,4-di-*tert*-pentylphenoxy) butyramido)-4,4-dimethyl-3-oxovaleranilide	402-260-4	54942-74-4	Aquatic Chronic 4	H413	—	H413			
616-025-00-0	valinamide	402-840-7	20108-78-5	Repr. 2 Eye Irrit. 2 Skin Sens. 1	H361f *** H319 H317	GHS08 Wng	H361f *** H319 H317			
616-026-00-6	thioacetamide	200-541-4	62-55-5	Carc. 1B Acute Tox. 4 * Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 3	H350 H302 H319 H315 H412	GHS08 GHS07 Dgr	H350 H302 H319 H315 H412			
616-027-00-1	tris(2-(2-hydroxyethoxy)ethyl)ammonium 3-acetoacetamido-4-methoxybenzenesulfonate	403-760-5	—	Skin Sens. 1	H317	GHS07 Wng	H317			
616-028-00-7	*N*-(4-(3-(4-cyanophenyl)ureido)-3-hydroxyphenyl)-2-(2,4-di-*tert*-pentylphenoxy)octanamide	403-790-9	108673-51-4	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
616-029-00-2	*N,N'*-ethylenebis(vinylsulfonylacetamide)	404-790-1	66710-66-5	Eye Dam. 1 Skin Sens. 1	H318 H317	GHS05 GHS07 Dgr	H318 H317			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-030-00-8	ethidimuron (ISO); 1-(5-ethylsulphonyl-1,3,4-thiadiazol-2-yl)-1,3-dimethylurea	250-010-6	30043-49-3	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
616-031-00-3	dimethachlor (ISO); 2-chloro-*N*-(2,6-dimethylphenyl)-*N*-(2-methoxyethyl)acetamide;	256-625-6	50563-36-5	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
616-032-00-9	diflufenican (ISO); N-(2,4-difluorophenyl)-2-[3-(trifluoromethyl)phenoxy]-3-pyridinecarboxamide	—	83164-33-4	Aquatic Chronic 3	H412	—	H412			
616-033-00-4	cyprofuram (ISO); *N*-(3-chlorophenyl)-*N*-(tetrahydro-2-oxo-3-furyl)cyclopropanecarboxamide	274-050-9	69581-33-5	Acute Tox. 3 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H301 H312 H400 H410	GHS06 GHS09 Dgr	H301 H312 H410			
616-034-00-X	pyracarbolid; (ISO); 3,4-dihydro-6-methyl-2*H*-pyran-5-carboxanilide	246-419-4	24691-76-7	Aquatic Chronic 3	H412	—	H412			
616-035-00-5	cymoxanil (ISO); 2-cyano-*N*-[(ethylamino)carbonyl]-2-(methoxyimino)acetamide	261-043-0	57966-95-7	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
616-036-00-0	2-chloracetamide	201-174-2	79-07-2	Repr. 2 Acute Tox. 3 * Skin Sens. 1	H361f *** H301 H317	GHS06 GHS08 Dgr	H361f *** H301 H317		Skin Sens. 1; H317: C ≥ 0,1 %	

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616-037-00-6	acetochlor (ISO); 2-chloro-*N*-(ethoxymethyl)-*N*-(2-ethyl-6-methylphenyl)acetamide	251-899-3	34256-82-1	Acute Tox. 4 * STOT SE 3 Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H332 H335 H315 H317 H400 H410	GHS07 GHS09 Wng	H332 H335 H315 H317 H410			
616-038-00-1	(4-aminophenyl)-*N*-methylmethylensulfonamide hydrochloride	406-010-5	88918-84-7	Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H318 H317 H411	GHS05 GHS07 GHS09 Dgr	H318 H317 H411			
616-039-00-7	3',5'-dichloro-4'-ethyl-2'-hydroxypalmitanilide	406-200-8	117827-06-2	Skin Sens. 1	H317	GHS07 Wng	H317			
616-040-00-2	potassium *N*-(4-toluenesulfonyl)-4-toluenesulfonamide	406-650-5	97888-41-0	Eye Dam. 1	H318	GHS05 Dgr	H318			
616-041-00-8	3',5'-dichloro-2-(2,4-di-*tert*-pentylphenoxy)-4'-ethyl-2'-hydroxyhexananilide	406-840-8	101664-25-9	Aquatic Chronic 4	H413	—	H413			
616-042-00-3	*N*-(2-(6-ethyl-7-(4-methylphenoxy)-1*H*-pyrazolo[1,5-b][1,2,4]triazol-2-yl)propyl)-2-octadecyloxybenzamide	407-070-5	142859-67-4	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
616-043-00-9	isoxaben (ISO); *N*-[3-(1-ethyl-1-methylpropyl)-1,2-oxazol-5-yl]-2,6-dimethoxybenzamide	407-190-8	82558-50-7	Aquatic Chronic 4	H413	—	H413			
616-044-00-4	*N*-(3,5-dichloro-4-ethyl-2-hydroxyphenyl)-2-(3-pentadecylphenoxy)-butanamide	402-510-2	—	Aquatic Chronic 2	H411	GHS09	H411			
616-045-00-X	2'-(4-chloro-3-cyano-5-formyl-2-thienylazo)-5'-diethylamino-2-methoxyacetanilide	405-190-2	122371-93-1	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			

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616-046-00-5	N-(2-(6-chloro-7-methylpyrazolo(1,5-b)-1,2,4-triazol-4-yl)propyl)-2-(2,4-di-*tert*-pentylphenoxy)octanamide	406-390-2	—	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-047-00-0	reaction mass of: 2,2',2",2"'-(ethylenedinitrilotetrakis-*N*,*N*-di(C_{16})alkylacetamide; 2,2',2",2"'-(ethylenedinitrilotetrakis-*N*,*N*-di(C_{18})alkylacetamide	406-640-0	—	Skin Sens. 1	H317	GHS07 Wng	H317			
616-048-00-6	3'-trifluoromethylisobutyranilide	406-740-4	1939-27-1	STOT RE 2 * Aquatic Chronic 2	H373 ** H411	GHS08 GHS09 Wng	H373 ** H411			
616-049-00-1	2-(2,4-bis(1,1-dimethylethyl)phenoxy)-*N*-(3,5-dichloro-4-ethyl-2-hydroxyphenyl)-hexanamide	408-150-2	99141-89-6	Aquatic Chronic 4	H413	—	H413			
616-050-00-7	lufenuron (ISO); *N*-[2,5-dichloro-4-(1,1,2,3,3,3-hexafluoropropoxy)-phenyl-aminocarbonyl]-2,6-difluorobenzamide	410-690-9	103055-07-8	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
616-051-00-2	reaction mass of: 2,4 -bis(*N*'-(4-methylphenyl)-ureido)-toluene; 2,6 -bis(*N*'-(4-methylphenyl)-ureido)-toluene	411-070-0	—	Aquatic Chronic 4	H413	—	H413			
616-052-00-8	formamide	200-842-0	75-12-7	Repr. 1B	H360-D ***	GHS08 Dgr	H360D ***			
616-053-00-3	*N*-methylacetamide	201-182-6	79-16-3	Repr. 1B	H360-D ***	GHS08 Dgr	H360D ***			
616-054-00-9	iprodione (ISO); 3-(3,5-dichlorophenyl)-2,4-dioxo-*N*-isopropylimidazolidine-1-carboxamide	253-178-9	36734-19-7	Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H400 H410	GHS08 GHS09 Wng	H351 H410			

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616-055-00-4	propyzamide (ISO); 3,5-dichloro-*N*-(1,1-dimethylprop-2-ynyl) benzamide	245-951-4	23950-58-5	Carc. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H400 H410	GHS08 GHS09 Wng	H351 H410			
616-056-00-X	*N*-methylformamide	204-624-6	123-39-7	Repr. 1B Acute Tox. 4 *	H360-D *** H312	GHS08 GHS07 Dg	H360D *** H312			
616-057-00-5	reaction mass of: *N*-[3-hydroxy-2-(2-methylacryloylaminomethoxy)propoxymethyl]-2-methylacrylamide; *N*-[2,3-bis-(2-methylacryloylaminomethoxy) propoxymethyl]-2-methylacrylamide; methacrylamide; 2-methyl-*N*-(2-methylacryloylamino-methoxymethyl)-acrylamide; *N*-(2,3-dihydroxypropoxymethyl)-2-methylacrylamide	412-790-8	—	Carc. 1B Muta. 2 STOT RE 2 *	H350 H341 H373 **	GHS08 Dgr	H350 H341 H373 **			
616-058-00-0	1,3-bis(3-methyl-2,5-dioxo-1*H*-pyrrolinyl-methyl)benzene	412-570-1	119462-56-5	STOT RE 2 * Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H373 ** H318 H317 H400 H410	GHS08 GHS05 GHS07 GHS09 Dgr	H373 ** H318 H317 H410			
616-059-00-6	4-((4-(diethylamino)-2-ethoxyphenyl) imino)-1,4-dihydro-1-oxo-*N*-propyl-2-naphthalenecarboxamide	412-650-6	121487-83-0	Aquatic Chronic 4	H413	—	H413			
616-060-00-1	Condensation product of: 3-(7-carboxyhept-1-yl)-6-hexyl-4-cyclohexene-1,2-dicarboxylic acid with polyamines (primarily amino-ethyl-piperazine and triethylenetetramine)	413-770-1	—	Acute Tox. 4 * Skin Corr. 1B Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H314 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H314 H317 H410			

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616-061-00-7	*N*,*N*'—1,6-hexanediylbis(*N*-(2,2,6,6-tetramethyl-piperidin-4-yl)-formamide	413-610-0	124172-53-8	Eye Irrit. 2 Aquatic Chronic 3	H319 H412	GHS07 Wng	H319 H412			
616-062-00-2	*N*-[3-[(2-acetyloxy)ethyl](phenyl-methyl) amino]-4-methoxyphenylacetamide	411-590-8	70693-57-1	Skin Corr. 1B Aquatic Chronic 3	H314 H412	GHS05 Dgr	H314 H412			
616-063-00-8	3-dodecyl-(1-(1,2,2,6,6-pentamethyl-4-piperidin)-yl)-2,5-pyrrolidindione	411-920-0	106917-30-0	Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1A Aquatic Acute 1 Aquatic Chronic 1	H331 H302 H373 ** H314 H400 H410	GHS06 GHS08 GHS05 GHS09 Dgr	H331 H302 H373 ** H314 H410			
616-064-00-3	*N*—*tert*-butyl-3-methylpicolinamide	406-720-5	32998-95-1	Aquatic Chronic 3	H412	—	H412			
616-065-00-9	3'-(3-acetyl-4-hydroxyphenyl)-1,1-diethylurea	411-970-3	79881-89-3	Acute Tox. 4 * STOT RE 2 *	H302 H373 **	GHS08 GHS07 Wng	H302 H373 **			
616-066-00-4	5,6,12,13-tetrachloroanthra(2,1,9-*def*:6,5,10-*d'e'f'*)diisoquinoline-1,3,8,10 (2*H*,9*H*)-tetrone	405-100-1	115662-06-1	Repr. 2	H361f ***	GHS08 Wng	H361f ***			
616-067-00-X	dodecyl 3-(2-(3-benzyl-4-ethoxy-2,5-dioxoimidazolidin-1-yl)-4,4-dimethyl-3-oxovaleramido)-4-chlorobenzoate	407-300-4	92683-20-0	Aquatic Chronic 4	H413	—	H413			
616-068-00-5	potassium 4-(11-methacrylamidoundecanamido)benzenesulfonate	406-500-9	174393-75-0	Skin Sens. 1	H317	GHS07 Wng	H317			
616-069-00-0	1-hydroxy-5-(2-methylpropyloxycarbonylamino)-*N*-(3-dodecyloxypropyl)-2-naphthoamide	406-210-2	110560-22-0	Aquatic Chronic 4	H413	—	H413			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-070-00-6	reaction mass of: 3,3'-dicyclohexyl-1,1'-methylenebis(4,1-phenylene)diurea; 3-cyclohexyl-1-(4-(4-(3-octadecylureido) benzyl)phenyl)urea; 3,3'-dioctadecyl-1,1'-methylenebis(4,1-phenylene)diurea	406-530-2	—	Aquatic Chronic 4	H413	—	H413			
616-071-00-1	reaction mass of: bis(*N*-cyclohexyl-*N*'-phenyleneureido)methylene; bis(*N*-octadecyl-*N*'-phenyleneureido)methylene; bis(*N*-dicyclohexyl-*N*'-phenyleneureido) methylene (1:2:1)	406-550-1	—	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
616-072-00-7	1-(2-deoxy-5-O-trityl-β-D-threopentofuranosyl)thymine	407-120-6	55612-11-8	Aquatic Chronic 4	H413	—	H413			
616-073-00-2	4'-ethoxy-2-benzimidazoleanilide	407-600-5	120187-29-3	Muta. 2 Aquatic Chronic 4	H341 H413	GHS08 Wng	H341 H413			
616-074-00-8	*N*-butyl-2-(4-morpholinylcarbonyl)benzamide	407-730-2	104958-67-0	Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 3	H319 H317 H412	GHS07 Wng	H319 H317 H412			
616-075-00-3	D, L-(*N*,*N*-diethyl-2-hydroxy-2-phenylacetamide)	408-120-9	65197-96-8	Acute Tox. 4 * Eye Dam. 1	H302 H318	GHS05 GHS07 Dgr	H302 H318			
616-076-00-9	tebufenozide (ISO); *N*-*tert*-butyl-*N*'-(4-ethylbenzoyl)-3,5-dimethylbenzohydrazide	412-850-3	112410-23-8	Aquatic Chronic 2	H411	GHS09	H411			

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616-077-00-4	reaction mass of: 2-(9-methyl-1,3,8,10-tetraoxo-2,3,9,10-tetrahydro-(1*H*,8*H*)-anthra[2,1,9-*def*: 6,5,10-*d'e'f'*]diisoquinolin-2-ylethansulfonic acid; potassium 2-(9-methyl-1,3,8,10-tetraoxo-2,3,9,10-tetrahydro-(1*H*,8*H*)-anthra[2,1,9-*def*: 6,5,10-*d'e'f'*]diisoquinolin-2-ylethansulfate	411-310-4	—	Eye Dam. 1	H318	GHS05 Dgr	H318			
616-078-00-X	2-[2,4-bis(1,1-dimethyl-ethyl)phenoxy]-*N*-(2-hydroxy-5-methyl-phenyl)hexanamide	411-330-3	104541-33-5	Aquatic Chronic 4	H413	—	H413			
616-079-00-5	1,6-hexanediyl-bis(2-(2-(1-ethylpentyl)-3-oxazolidinyl)ethyl)carbamate	411-700-4	140921-24-0	Skin Sens. 1	H317	GHS07 Wng	H317			
616-080-00-0	4-(2-((3-ethyl-4-methyl-2-oxo-pyrrolin-1-yl)carboxamido)ethyl)benzenesulfonamide)	411-850-0	119018-29-0	Aquatic Chronic 3	H412	—	H412			
616-081-00-6	5-bromo-8-naphtholactam	413-480-5	24856-00-6	Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H317 H400 H410	GHS07 GHS09 Wng	H302 H317 H410			
616-082-00-1	*N*-(5-chloro-3-((4-(diethylamino)-2-methylphenyl)imino-4-methyl-6-oxo-1,4-cyclohexadien-1-yl)benzamide	413-200-1	129604-78-0	Skin Sens. 1	H317	GHS07 Wng	H317			
616-083-00-7	[2-[(4-nitrophenyl)amino]ethyl]urea	410-700-1	27080-42-8	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
616-084-00-2	2,4-bis[*N*'-(4-methylphenyl)ureido]toluene	411-790-5	—	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-085-00-8	3-(2,4-dichlorophenyl)-6-fluoro-quinazoline-2,4(1*H*,3*H*)-dione	412-190-6	168900-02-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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616-086-00-3	2-acetylamino-6-chloro-4-[(4-diethylamino) 2-methylphenyl-imino]-5-methyl-1-oxo-2,5-cyclohexadiene	412-250-1	102387-48-4	Aquatic Chronic 4	H413	—	H413			
616-087-00-9	reaction mass of: 7,9,9-trimethyl-3,14-dioxa-4,13-dioxo-5,12-diazahexadecane-1,16-diyl-prop-2-enoate; 7,7,9-trimethyl-3,14-dioxa-4,13-dioxo-5,12-diazahexadecan-1,16-diyl-prop-2-enoate	412-260-6	52658-19-2	Eye Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H319 H317 H411	GHS07 GHS09 Wng	H319 H317 H411			
616-088-00-4	2-aminosulfonyl-*N*,*N*-dimethylnicotinamide	413-440-7	112006-75-4	Skin Sens. 1 Aquatic Chronic 3	H317 H412	GHS07 Wng	H317 H412			
616-089-00-X	5-(2,4-dioxo-1,2,3,4-tetrahydropyrimidine)-3-fluoro-2-hydroxymethyltetrahydrofuran	415-360-8	41107-56-6	Muta. 2	H341	GHS08 Wng	H341			
616-090-00-5	1-(1,4-benzodioxan-2-ylcarbonyl)piperazine hydrochloride	415-660-9	70918-74-0	Acute Tox. 3 * Acute Tox. 3 * Acute Tox. 3 * STOT RE 2 * Aquatic Chronic 2	H331 H311 H301 H373 ** H411	GHS06 GHS08 GHS09 Dgr	H331 H311 H301 H373 ** H411			
616-091-00-0	1,3,5-tris-[(2*S* and 2*R*)-2,3-epoxypropyl]-1,3,5-triazine-2,4,6-(1*H*,3*H*,5*H*)-trione	423-400-0	59653-74-6	Muta. 1B Acute Tox. 3 * Acute Tox. 4 * STOT RE 2 * Eye Dam. 1 Skin Sens. 1	H340 H331 H302 H373 ** H318 H317	GHS06 GHS08 GHS05 Dgr	H340 H331 H302 H373 ** H318 H317			
616-092-00-6	Polymeric reaction product of bicyclo[2.2.1] hepta-2,5-diene, ethene, 1,4-hexadiene, 1-propene with *N*,*N*-di-2-propenylformamide	404-035-6	—	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
616-093-00-1	Reaction products of: aniline-terephthalaldehyde-*o*-toluidine condensate with maleic anhydride	406-620-1	129217-90-9	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			

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616-094-00-7	3,3'-dicyclohexyl-1,1'-methylenebis(4,1-phenylene)diurea	406-370-3	58890-25-8	Skin Sens. 1 Aquatic Chronic 4	H317 H413	GHS07 Wng	H317 H413			
616-095-00-2	3,3'-dioctadecyl-1,1'-methylenebis(4,1-phenylene)diurea	406-690-3	43136-14-7	Aquatic Chronic 4	H413	—	H413			
616-096-00-8	*N*-(3-hexadecyloxy-2-hydroxyprop-1-yl)-*N*-(2-hydroxyethyl)palmitamide	408-110-4	110483-07-3	Aquatic Chronic 4	H413	—	H413			
616-097-00-3	*N*,*N*'-1,4-phenylenebis(2-((2-methoxy-4-nitrophenyl)azo)-3-oxobutanamide	411-840-6	83372-55-8	Aquatic Chronic 4	H413	—	H413			
616-098-00-9	1-[4-chloro-3-((2,2,3,3,3-pentafluoropropoxy)methyl)phenyl]-5-phenyl-1*H*-1,2,4-triazole-3-carboxamide	411-750-7	119126-15-7	Aquatic Chronic 2	H411	GHS09	H411			
616-099-00-4	2-[4-[(4-hydroxyphenyl)sulfonyl]phenoxy]-4,4-dimethyl-*N*-[5-[(methylsulfonyl)amino]-2-[4-(1,1,3,3-tetramethylbutyl)phenoxy]phenyl]-3-oxopentanamide	414-170-2	135937-20-1	Aquatic Chronic 4	H413	—	H413			
616-100-00-8	1,3-dimethyl-1,3-bis(trimethylsilyl)urea	414-180-7	10218-17-4	Acute Tox. 4 * Skin Irrit. 2	H302 H315	GHS07 Wng	H302 H315			
616-101-00-3	(*S*)-*N*-*tert*-butyl-1,2,3,4-tetrahydro-3-isoquinolinecarboxamide	414-600-9	149182-72-9	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
616-102-00-9	reaction mass of: α-[3-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyl]-ω-[3-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyloxy]-poly-(oxyethylene-co-oxypropylene); 1,2-(or 1,3-)bis[α-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyl)-ω-oxy-poly(oxyethylene-co-oxypropylene)]-3-(or 2-)propanol; 1,2,3-tris[α-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyl)-ω-oxy-poly-(oxyethylene-co-oxypropylene)]propane]	415-870-0	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-103-00-4	(*S*,*S*)-*trans*-4-(acetylamino)-5,6-dihydro-6-methyl-7,7-dioxo-4*H*-thieno[2,3-b]thiopyran-2-sulfonamide	415-030-3	120298-38-6	Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H317 H400 H410	GHS07 GHS09 Wng	H317 H410			
616-104-00-X	benalaxyl (ISO); methyl *N*-(2,6-dimethylphenyl)-*N*-(phenylacetyl)-DL-alaninate	275-728-7	71626-11-4	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-105-00-5	chlorotoluron (ISO); 3-(3-chloro-*p*-tolyl)-1,1-dimethylurea	239-592-2	15545-48-9	Carc. 2 Repr. 2 Aquatic Acute 1 Aquatic Chronic 1	H351 H361d *** H400 H410	GHS08 GHS09 Wng	H351 H361d *** H410			
616-106-00-0	phenmedipham (ISO); methyl 3-(3-methylcarbaniloyloxy)carbanilate	237-199-0	13684-63-4	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-108-00-1	iodosulfuron-methyl-sodium; sodium ({[5-iodo-2-(methoxycarbonyl)phenyl]sulfonyl}carbamoyl)(4-methoxy-6-methyl-1,3,5-triazin-2-yl)azanide	—	144550-36-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-109-00-7	sulfosulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-(2-ethylsulfonylimidazo[1,2-a]pyridin-3-yl)sulfonylurea	—	141776-32-1	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-110-00-2	cyclanilide (ISO); 1-(2,4-dichloroanilinocarbonyl)cyclopropanecarboxylic acid	419-150-7	113136-77-9	Acute Tox. 4 * Aquatic Chronic 2	H302 H411	GHS07 GHS09 Wng	H302 H411			
616-111-00-8	fenhexamid (ISO); *N*-(2,3-dichlor-4-hydroxyphenyl)-1-methylcyclohexancarboxamid	422-530-5	126833-17-8	Aquatic Chronic 2	H411	GHS09	H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-112-00-3	oxasulfuron (ISO); oxetan-3-yl 2-[(4,6-dimethylpyrimidin-2-yl)-carbamoylsulfamoyl]benzoate	—	144651-06-9	STOT RE 2 * Aquatic Acute 1 Aquatic Chronic 1	H373 ** H400 H410	GHS08 GHS09 Wng	H373 ** H410			
616-113-00-9	desmedipham (ISO); ethyl 3-phenylcarbamoyloxyphenylcarbamate	237-198-5	13684-56-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410		M=10	
616-114-00-4	dodecanamide, *N,N'*-(9,9',10,10'-tetrahydro-9,9',10,10'-tetraoxo(1,1'-bianthracene)-4,4'-diyl)bis-	418-010-2	136897-58-0	Aquatic Chronic 4	H413	—	H413			
616-115-00-X	*N*-(3-acetyl-2-hydroxyphenyl)-4-(4-phenylbutoxy)benzamide	416-150-9	136450-06-1	Aquatic Chronic 4	H413	—	H413			
616-116-00-5	*N*-(4-dimethylaminopyridinium)-3-methoxy-4-(1-methyl-5-nitroindol-3-ylmethyl)-*N*-(*o*-tolylsulfonyl)benzamidate	416-790-9	143052-96-4	Aquatic Chronic 4	H413	—	H413			
616-117-00-0	*N*-[2-(3-acetyl-5-nitrothiophen-2-ylazo)-5-diethylaminophenyl]acetamide	416-860-9	777891-21-1	Repr. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H361f *** H317 H400 H410	GHS08 GHS09 Wng	H361f *** H317 H410			
616-118-00-6	*N*-(2',6'-dimethylphenyl)-2-piperidinecarboxamide hydrochloride	417-950-0	65797-42-4	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
616-119-00-1	2-(1-butyl-3,5-dioxo-2-phenyl-(1,2,4)-triazolidin-4-yl)-4,4-dimethyl-3-oxo-*N*-(2-methoxy-5-(2-(dodecyl-1-sulfonyl))propionylamino)-phenyl)-pentanamide	418-060-5	118020-93-2	Aquatic Chronic 4	H413	—	H413			
616-120-00-7	reaction mass of: *N*-(3-dimethylamino-4-methyl-phenyl)-benzamide; *N*-(3-dimethylamino-2-methyl-phenyl)-benzamide; *N*-(3-dimethylamino-3-methyl-phenyl)-benzamide	420-600-1	—	STOT RE 2 * Aquatic Chronic 2	H373 ** H411	GHS08 GHS09 Wng	H373 ** H411			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-121-00-2	2,4-dihydroxy-*N*-(2-methoxyphenyl)benzamide	419-090-1	129205-19-2	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
616-123-00-3	*N*-[3-[[4-(diethylamino)-2-methylphenyl]imino]-6-oxo-1,4-cyclohexadienyl]acetamide	414-740-0	96141-86-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
616-124-00-9	lithium bis(trifluoromethylsulfonyl)imide	415-300-0	90076-65-6	Acute Tox. 3 * Acute Tox. 3 * Skin Corr. 1B Aquatic Chronic 3	H311 H301 H314 H412	GHS06 GHS05 Dgr	H311 H301 H314 H412			
616-125-00-4	3-cyano-*N*-(1,1-dimethylethyl)androsta-3,5-diene-17-β-carboxamide	415-730-9	151338-11-3	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	410			
616-127-00-5	reaction mass of: *N*,*N*'-Ethane-1,2-diylbis(decanamide); 12-Hydroxy-*N*-[2-[1-oxydecyl)amino]ethyl]octadecanamide; *N*,*N*'-Ethane-1,2-diylbis(12-hydroxyoctadecanamide)	430-050-2	—	Skin Sens. 1 Aquatic Chronic 2	H317 H411	GHS07 GHS09 Wng	H317 H411			
616-128-00-0	*N*-(2-(1-allyl-4,5-dicyanoimidazol-2-ylazo)-5-(dipropylamino)phenyl)-acetamide	417-530-7	123590-00-1	Aquatic Chronic 4	H413	—	H413			
616-129-00-6	*N*,*N*'-bis(2,2,6,6-tetramethyl-4-piperidyl)isophthalamide	419-710-0	42774-15-2	Acute Tox. 4 * Eye Irrit. 2	H302 H319	GHS07 Wng	H302 H319			
616-130-00-1	*N*-(3-(2-(4,4-dimethyl-2,5-dioxo-imidazolin-1-yl)-4,4-dimethyl-3-oxo-pentanoylamino)-4-methoxy-phenyl)-octadecanamide	421-780-2	150919-56-5	Aquatic Chronic 4	H413	—	H413			
616-132-00-2	*N*-[4-(4-cyano-2-furfurylidene-2,5-dihydro-5-oxo-3-furyl)phenyl]butane-1-sulfonamide	423-250-6	130016-98-7	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
616-133-00-8	*N*-cyclohexyl-*S*,*S*-dioxobenzo[*b*]thiophene-2-carboxamide	423-990-1	149118-66-1	Acute Tox. 4 * Eye Dam. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H318 H400 H410	GHS05 GHS07 GHS09 Dgr	H302 H318 H410			
616-134-00-3	3,3'-bis(dioctyloxyphosphinothioylthio)-*N*,*N*'-oxybis(methylene)dipropionamide	401-820-5	793710-14-2	Aquatic Chronic 3	H412	—	H412			
616-135-00-9	(3*S*,4a*S*,8a*S*)-2-[(2*R*,3*S*)-3-amino-2-hydroxy-4-phenylbutyl]-*N*-tert-butyldecahydroisoquinoline-3-carboxamide	430-230-0	136522-17-3	Acute Tox. 4 * Aquatic Chronic 3	H302 H412	GHS07 Wng	H302 H412			
616-142-00-7	1,3-Bis(vinylsulfonylacetamido)propane	428-350-3	93629-90-4	Muta. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 3	H341 H318 H317 H412	GHS08 GHS05 GHS07 Dgr	H341 H318 H317 H412			
616-143-00-2	*N*,*N*'-dihexadecyl-*N*,*N*'-bis(2-hydroxyethyl) propanediamide	422-560-9	149591-38-8	Repr. 2 Eye Irrit. 2 Aquatic Chronic 4	H361f *** H319 H413	GHS08 Wng	H361f *** H319 H413			
617-001-00-2	di-*tert*-butyl peroxide	203-733-6	110-05-4	Org. Perox. E Flam. Liq. 2	H242 H225	GHS02 Dgr	H242 H225			
617-002-00-8	α, α-dimethylbenzyl hydroperoxide; cumene hydroperoxide	201-254-7	80-15-9	Org. Perox. E Acute Tox. 3 * Acute Tox. 4 * Acute Tox. 4 * STOT RE 2 * Skin Corr. 1B Aquatic Chronic 2	H242 H331 H312 H302 H373 ** H314 H411	GHS02 GHS06 GHS08 GHS05 GHS09 Dgr	H242 H331 H312 H302 H373 ** H314 H411		Skin Corr. 1B; H314: C ≥ 10 % Skin Irrit. 2; H315: 3 % ≤ C < 10 % Eye Dam. 1; H318: 3 % ≤ C < 10 % Eye Irrit. 2; H319: 1 % ≤ C < 3 % STOT SE 3; H335: C < 10 %	

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
617-003-00-3	dilauroyl peroxide	203-326-3	105-74-8	Org. Perox. D	H242	GHS02 Dgr	H242			
617-004-00-9	1,2,3,4-tetrahydro-1-naphthyl hydroperoxide	212-230-0	771-29-9	Org. Perox. D Acute Tox. 4 * Skin Corr. 1B Aquatic Acute 1 Aquatic Chronic 1	H242 H302 H314 H400 H410	GHS02 GHS05 GHS07 GHS09 Dgr	H242 H302 H314 H410		STOT SE 3; H335: C ≥ 5 %	
617-006-00-X	bis(α, α-dimethylbenzyl) peroxide	201-279-3	80-43-3	Org. Perox. F Eye Irrit. 2 Skin Irrit. 2 Aquatic Chronic 2	H242 H319 H315 H411	GHS02 GHS07 GHS09 Wng	H242 H319 H315 H411			
617-007-00-5	*tert*-butyl α, α-dimethylbenzyl peroxide	222-389-8	3457-61-2	Org. Perox. E Skin Irrit. 2 Aquatic Chronic 2	H242 H315 H411	GHS02 GHS07 GHS09 Wng	H242 H315 H411			
617-008-00-0	dibenzoyl peroxide; benzoyl peroxide	202-327-6	94-36-0	Org. Perox. B Eye Irrit. 2 Skin Sens. 1	H241 H319 H317	GHS01 GHS02 GHS07 Wng	H241 H319 H317			T
617-010-00-1	1-hydroperoxycyclohexyl 1-hydroxycyclohexyl peroxide; [1] 1,1'-dioxybiscyclohexan-1-ol; [2] cyclohexylidene hydroperoxide; [3] cyclohexanone, peroxide [4] [> 91 % solution]	201-091-1 [1] 219-306-2 [2] 220-279-4 [3] 235-527-7 [4]	78-18-2 [1] 2407-94-5 [2] 2699-11-8 [3] 12262-58-7 [4]	Org. Perox. A Acute Tox. 4 * Skin Corr. 1B	H240 H302 H314	GHS01 GHS05 GHS07 Dgr	H240 H302 H314		STOT SE 3; H335: C ≥ 5 %	C T

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
617-010-01-9	1-hydroperoxycyclohexyl 1-hydroxycyclohexyl peroxide; [1] 1,1'-dioxybiscyclohexan-1-ol; [2] cyclohexylidene hydroperoxide; [3] cyclohexanone, peroxide [4] [≤ 91 % solution]	201-091-1 [1] 219-306-2 [2] 220-279-4 [3] 235-527-7 [4]	78-18-2 [1] 2407-94-5 [2] 2699-11-8 [3] 12262-58-7 [4]	Org. Perox. C Acute Tox. 4 * Skin Corr. 1B	H242 H302 H314	GHS02 GHS05 GHS07 Dgr	H242 H302 H314		STOT SE 3; H335: C ≥ 5 %	C T
617-012-00-2	8-*p*-menthyl hydroperoxide; *p*-menthane hydroperoxide	201-281-4	80-47-7	Org. Perox. D Skin Corr. 1B Acute Tox. 4 *	H242 H314 H332	GHS02 GHS05 GHS07 Dgr	H242 H314 H332		STOT SE 3; H335: C ≥ 5 %	
617-013-00-8	*O*,*O*-*tert*-butyl *O*-docosyl monoperoxyoxalate	404-300-6	116753-76-5	Org. Perox. C **** Aquatic Acute 1 Aquatic Chronic 1	H242 H400 H410	GHS02 GHS09 Dgr	H242 H410			
617-014-00-3	6-(nonylamino)-6-oxo-peroxyhexanoic acid	406-680-9	104788-63-8	Org. Perox. C **** Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1	H242 H318 H317 H400	GHS02 GHS05 GHS07 GHS09 Dgr	H242 H318 H317 H400			
617-015-00-9	bis(4-methylbenzoyl)peroxide	407-950-9	895-85-2	Org. Perox. B **** Aquatic Acute 1 Aquatic Chronic 1	H241 H400 H410	GHS01 GHS02 GHS09 Dgr	H241 H410			
617-016-00-4	3-hydroxy-1,1-dimethylbutyl 2-ethyl-2-methylheptaneperoxoate	413-910-1	—	Org. Perox. C **** Flam. Liq. 3 Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H242 H226 H315 H400 H410	GHS02 GHS07 GHS09 Dgr	H242 H226 H315 H410			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
617-017-00-X	reaction mass of: 2,2'-bis(*tert*-pentylperoxy)-*p*-diisopropylbenzene; 2,2'-bis(*tert*-pentylperoxy)-*m*-diisopropylbenzene	412-140-3	32144-25-5	Org. Perox. D **** Aquatic Chronic 4	H242 H413	GHS02 Dgr	H242 H413			T
617-018-00-5	reaction mass of: 1-methyl-1-(3-(1-methylethyl)phenyl)ethyl-1-methyl-1-phenylethylperoxide, 63 % by weight; 1-methyl-1-(4-(1-methylethyl)phenyl)ethyl-1-methyl-1-phenylethylperoxide, 31 % by weight	410-840-3	71566-50-2	Org. Perox. C **** Aquatic Chronic 2	H242 H411	GHS02 GHS09 Dgr	H242 H411			T
617-019-00-0	6-(phthalimido)peroxyhexanoic acid	410-850-8	128275-31-0	Org. Perox. D Eye Dam. 1 Aquatic Acute 1	H242 H318 H400	GHS02 GHS05 GHS09 DgDgr	H242 H318 H400			T
617-020-00-6	1,3-di(prop-2,2-diyl)benzene bis(neodecanoylperoxide)	420-060-5	117663-11-3	Flam. Liq. 3 Org. Perox. D **** Aquatic Chronic 2	H226 H242 H411	GHS02 GHS09 Dgr	H226 H242 H411			
647-001-00-8	glucosidase, β-	232-589-7	9001-22-3	Resp. Sens. 1	H334	GHS08 Dgr	H334			
647-002-00-3	cellulase	232-734-4	9012-54-8	Resp. Sens. 1	H334	GHS08 Dgr	H334			
647-003-00-9	cellobiohydrolase, exo-	253-465-9	37329-65-0	Resp. Sens. 1	H334	GHS08 Dgr	H334			
647-004-00-4	cellulases with the exception of those specified elsewhere in this Annex	—	—	Resp. Sens. 1	H334	GHS08 Dgr	H334			A
647-005-00-X	bromelain, juice	232-572-4	9001-00-7	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			

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647-006-00-5	ficin	232-599-1	9001-33-6	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-007-00-0	papain	232-627-2	9001-73-4	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-008-00-6	pepsin A	232-629-3	9001-75-6	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-009-00-1	rennin	232-645-0	9001-98-3	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-010-00-7	trypsin	232-650-8	9002-07-7	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-011-00-2	chymotrypsin	232-671-2	9004-07-3	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-012-00-8	subtilisin	232-752-2	9014-01-1	STOT SE 3 Skin Irrit. 2 Eye Dam. 1 Resp. Sens. 1	H335 H315 H318 H334	GHS08 GHS05 GHS07 Dgr	H335 H315 H318 H334			

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
647-013-00-3	proteinase, microbial neutral	232-966-6	9068-59-1	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-014-00-9	proteases with the exception of those specified elsewhere in this Annex	—	—	Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Resp. Sens. 1	H319 H335 H315 H334	GHS08 GHS07 Dgr	H319 H335 H315 H334			
647-015-00-4	amylase, α-	232-565-6	9000-90-2	Resp. Sens. 1	H334	GHS08 Dgr	H334			
647-016-00-X	amylases with the exception of those specified elsewhere in this Annex	—	—	Resp. Sens. 1	H334	GHS08 Dgr	H334			
648-001-00-0	Distillates (coal tar), benzole fraction; Light Oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists of hydrocarbons having carbon numbers primarily in the range of C_4 to C_{10} and distilling in the approximate range of 80 °C to 160 °C (175 °F to 320 °F).]	283-482-7	84650-02-2	Carc. 1B	H350	GHS08 Dgr	H350			H
648-002-00-6	Tar oils, brown-coal; Light Oil; [The distillate from lignite tar boiling in the range of approximately 80 °C to 250 °C (176 °F to 482 °F). Composed primarily of aliphatic and aromatic hydrocarbons and monobasic phenols.]	302-674-4	94114-40-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-003-00-1	Benzol forerunnings (coal); Light Oil Redistillate, low boiling; [The distillate from coke oven light oil having an approximate distillation range below 100 °C (212 °F). Composed primarily of C_4 to C_6 aliphatic hydrocarbons.]	266-023-5	65996-88-5	Carc. 1B	H350	GHS08 Dgr	H350			H J

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648-004-00-7	Distillates (coal tar), benzole fraction, BTX-rich; Light Oil Redistillate, low boiling; [A residue from the distillation of crude benzole to remove benzole fronts. Composed primarily of benzene, toluene and xylenes boiling in the range of approximately 75 °C to 200 °C (167 °F to 392 °F).]	309-984-9	101896-26-8	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-005-00-2	Aromatic hydrocarbons, C_{6-10}, C_8-rich; Light Oil Redistillate, low boiling	292-697-5	90989-41-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-006-00-8	Solvent naphtha (coal), light; Light Oil Redistillate, low boiling	287-498-5	85536-17-0	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-007-00-3	Solvent naphtha (coal), xylene-styrene cut; Light Oil Redistillate, intermediate boiling	287-502-5	85536-20-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-008-00-9	Solvent naphtha (coal), coumarone-styrene contg.; Light Oil Redistillate, intermediate boiling	287-500-4	85536-19-2	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-009-00-4	Naphtha (coal), distn. residues; Light Oil Redistillate, high boiling; [The residue remaining from the distillation of recovered naphtha. Composed primarily of naphthalene and condensation products of indene and styrene.]	292-636-2	90641-12-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-010-00-X	Aromatic hydrocarbons, C_8; Light Oil Redistillate, high boiling	292-694-9	90989-38-1	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-012-00-0	Aromatic hydrocarbons, C_{8-9}, hydrocarbon resin polymn. by-product; Light Oil Redistillate, high boiling; [A complex combination of hydrocarbons obtained from the evaporation of solvent under vacuum from polymerized hydrocarbon resin. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_9 and boiling in the range of approximately 120 °C to 215 °C (248 °F to 419 °F).]	295-281-1	91995-20-9	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-013-00-6	Aromatic hydrocarbons, C_{9-12}, benzene distn.; Light Oil Redistillate, high boiling	295-551-9	92062-36-7	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-014-00-1	Extract residues (coal), benzole fraction alk., acid ext.; Light Oil Extract Residues, low boiling; [The redistillate from the distillate, freed of tar acids and tar bases, from bituminous coal high temperature tar boiling in the approximate range of 90 °C to 160 °C (194 °F to 320 °F). It consists predominantly of benzene, toluene and xylenes.]	295-323-9	91995-61-8	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-015-00-7	Extract residues (coal tar), benzole fraction alk., acid ext.; Light Oil Extract Residues, low boiling; [A complex combination of hydrocarbons obtained by the redistillation of the distillate of high temperature coal tar (tar acid and tar base free). It consists predominantly of unsubstituted and substituted mononuclear aromatic hydrocarbons boiling in the range of 85 °C-195 °C (185 °F-383 °F).]	309-868-8	101316-63-6	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-016-00-2	Extract residues (coal), benzole fraction acid; Light Oil Extract Residues, low boiling; [An acid sludge by-product of the sulphuric acid refining of crude high temperature coal. Composed primarily of sulfuric acid and organic compounds.]	298-725-2	93821-38-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-017-00-8	Extract residues (coal), light oil alk., distn. overheads; Light Oil Extract Residues, low boiling; [The first fraction from the distillation of aromatic hydrocarbons, coumarone, naphthalene and indene rich prefactionator bottoms or washed carbolic oil boiling substantially below 145 °C (293 °F). Composed primarily of C_7 and C_8 aliphatic and aromatic hydrocarbons.]	292-625-2	90641-02-4	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-018-00-3	Extract residues (coal), light oil alk., acid ext., indene fraction; Light Oil Extract Residues, intermediate boiling	309-867-2	101316-62-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-019-00-9	Extract residues (coal), light oil alk., indene naphtha fraction; Light Oil Extract Residues, high boiling; [The distillate from aromatic hydrocarbons, coumarone, naphthalene and indene rich prefractionator bottoms or washed carbolic oils, having an approximate boiling range of 155 °C to 180 °C (311 °F to 356 °F). Composed primarily of indene, indan and trimethylbenzenes.]	292-626-8	90641-03-5	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-020-00-4	Solvent naphtha (coal); Light Oil Extract Residues, high boiling; [The distillate from either high temperature coal tar, coke oven light oil, or coal tar oil alkaline extract residue having an approximate distillation range of 130 °C to 210 °C (266 °F to 410 °F) Composed primarily of indene and other polycyclic ring systems containing a single aromatic ring. May contain phenolic compounds and aromatic nitrogen bases.]	266-013-0	65996-79-4	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-021-00-X	Distillates (coal tar), light oils, neutral fraction; Light Oil Extract Residues, high boiling; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of alkyl-substituted one ring aromatic hydrocarbons boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F). May also include unsaturated hydrocarbons such as indene and coumarone.]	309-971-8	101794-90-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-022-00-5	Distillates (coal tar), light oils, acid exts.; Light Oil Extract Residues, high boiling; [This oil is a complex mixture of aromatic hydrocarbons, primarily indene, naphthalene, coumarone, phenol, and *o*-, *m*- and *p*-cresol and boiling in the range of 140 °C to 215 °C (284 °F to 419 °F).]	292-609-5	90640-87-2	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-023-00-0	Distillates (coal tar), light oils; Carbolic Oil; [A complex combination of hydrocarbons obtained by distillation of coal tar. It consists of aromatic and other hydrocarbons, phenolic compounds and aromatic nitrogen compounds and distills at the approximate range of 150 °C to 210 °C (302 °F to 410 °F).]	283-483-2	84650-03-3	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-024-00-6	Tar oils, coal; Carbolic Oil; [The distillate from high temperature coal tar having an approximate distillation range of 130 °C to 250 °C (266 °F to 410 °F). Composed primarily of naphthalene, alkylnaphthalenes, phenolic compounds, and aromatic nitrogen bases.]	266-016-7	65996-82-9	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-026-00-7	Extract residues (coal), light oil alk., acid ext.; Carbolic Oil Extract Residue; [The oil resulting from the acid washing of alkali-washed carbolic oil to remove the minor amounts of basic compounds (tar bases). Composed primarily of indene, indan and alkylbenzenes.]	292-624-7	90641-01-3	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-027-00-2	Extract residues (coal), tar oil alk.; Carbolic Oil Extract Residue; [The residue obtained from coal tar oil by an alkaline wash such as aqueous sodium hydroxide after the removal of crude coal tar acids. Composed primarily of naphthalenes and aromatic nitrogen bases.]	266-021-4	65996-87-4	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-028-00-8	Extract oils (coal), light oil; Acid Extract; [The aqueous extract produced by an acidic wash of alkali-washed carbolic oil. Composed primarily of acid salts of various aromatic nitrogen bases including pyridine, quinoline and their alkyl derivatives.]	292-622-6	90640-99-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-029-00-3	Pyridine, alkyl derivs.; Crude Tar Bases; [The complex combination of polyalkylated pyridines derived from coal tar distillation or as high-boiling distillates approximately above 150 °C (302 °F) from the reaction of ammonia with acetaldehyde, formaldehyde or paraformaldehyde.]	269-929-9	68391-11-7	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-030-00-9	Tar bases, coal, picoline fraction; Distillate Bases; [Pyridine bases boiling in the range of approximately 125 °C to 160 °C (257 °F 320 °F) obtained by distillation of neutralized acid extract of the base-containing tar fraction obtained by the distillation of bituminous coal tars. Composed chiefly of lutidines and picolines.]	295-548-2	92062-33-4	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-031-00-4	Tar bases, coal, lutidine fraction; Distillate Bases	293-766-2	91082-52-9	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-032-00-X	Extract oils (coal), tar base, collidine fraction; Distillate Bases; [The extract produced by the acidic extraction of bases from crude coal tar aromatic oils, neutralization, and distillation of the bases. Composed primarily of collidines, aniline, toluidines, lutidines, xylidines.]	273-077-3	68937-63-3	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-033-00-5	Tar bases, coal, collidine fraction; Distillate Bases; [The distillation fraction boiling in the range of approximately 181 °C to 186 °C (356 °F to 367 °F) from the crude bases obtained from the neutralized, acid-extracted base-containing tar fractions obtained by the distillation of bituminous coal tar. It contains chiefly aniline and collidines.]	295-543-5	92062-28-7	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-034-00-0	Tar bases, coal, aniline fraction; Distillate Bases; [The distillation fraction boiling in the range of approximately 180 °C to 200 °C (356 °F to 392 °F) from the crude bases obtained by dephenolating and debasing the carbolated oil from the distillation of coal tar. It contains chiefly aniline, collidines, lutidines and toluidines.]	295-541-4	92062-27-6	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-035-00-6	Tar bases, coal, toluidine fraction; Distillate Bases	293-767-8	91082-53-0	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-036-00-1	Distillates (petroleum), alkene-alkyne manuf. pyrolysis oil, mixed with high-temp. coal tar, indene fraction; Redistillates; [A complex combination of hydrocarbons obtained as a redistillate from the fractional distillation of bituminous coal high temperature tar and residual oils that are obtained by the pyrolytic production of alkenes and alkynes from petroleum products or natural gas. It consists predominantly of indene and boils in a range of approximately 160 °C to 190 °C (320 °F to 374 °F).]	295-292-1	91995-31-2	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-037-00-7	Distillates (coal), coal tar-residual pyrolysis oils, naphthalene oils; Redistillates; [The redistillate obtained from the fractional distillation of bituminous coal high temperature tar and pyrolysis residual oils and boiling in the range of approximately 190 °C to 270 °C (374 °F to 518 °F). Composed primarily of substituted dinuclear aromatics.]	295-295-8	91995-35-6	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-038-00-2	Extract oils (coal), coal tar-residual pyrolysis oils, naphthalene oil, redistillate; Redistillates; [The redistillate from the fractional distillation of dephenolated and debased methylnaphthalene oil obtained from bituminous coal high temperature tar and pyrolysis residual oils boiling in the approximate range of 220 °C to 230 °C (428 °F to 446 °F). It consists predominantly of unsubstituted and substituted dinuclear aromatic hydrocarbons.]	295-329-1	91995-66-3	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-039-00-8	Extract oils (coal), coal tar-residual pyrolysis oils, naphthalene oils; Redistillates; [A neutral oil obtained by debasing and dephenolating the oil obtained from the distillation of high temperature tar and pyrolysis residuel oils which has a boiling range of 225 °C to 255 °C (437 °F to 491 °F). Composed primarily of substituted dinuclear aromatic hydrocarbons.]	310-170-0	122070-79-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-040-00-3	Extract oils (coal), coal tar residual pyrolysis oils, naphthalene oil, distn. residues; Redistillates; [Residue from the distillation of dephenolated and debased methylnaphthalene oil (from bituminous coal tar and pyrolysis residual oils) with a boiling range of 240 °C to 260 °C (464 °F to 500 °F). Composed primarily of substituted dinuclear aromatic and heterocyclic hydrocarbons.]	310-171-6	122070-80-8	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-041-00-9	Absorption oils, bicyclo arom. and heterocyclic hydrocarbon fraction; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained as a redistillate from the distillation of wash oil. It consists predominantly of 2-ringed aromatic and heterocyclic hydrocarbons boiling in the range of approximately 260 °C to 290 °C (500 °F to 554 °F).]	309-851-5	101316-45-4	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-042-00-4	Distillates (coal tar), upper, fluorene-rich; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained by the crystallization of tar oil. It consists af aromatic and polycyclic hydrocarbons primarily fluorene and some acenaphthene.]	284-900-0	84989-11-7	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-043-00-X	Creosote oil, acenaphthene fraction, acenaphthene-free; Wash Oil Redistillate; [The oil remaining after removal by a crystallization process of acenaphthene from acenaphthene oil from coal tar. Composed primarily of naphthalene and alkylnaphthalenes.]	292-606-9	90640-85-0	Carc. 1B	H350	GHS08 Dgr	H350			H
648-044-00-5	Distillates (coal tar), heavy oils; Heavy Anthracene Oil; [Distillate from the fractional distillation of coal tar of bituminous coal, with boiling range of 240 °C to 400 °C (464 °F to 752 °F). Composed primarily of tri- and polynuclear hydrocarbons and heterocyclic compounds.]	292-607-4	90640-86-1	Carc. 1B	H350	GHS08 Dgr	H350			H
648-045-00-0	Distillates (coal tar), upper; Heavy Anthracene Oil; [The distillate from coal tar having an approximate distillation range of 220 °C to 450 °C (428 °F to 842 °F). Composed primarily of three to four membered condensed ring aromatic hydrocarbons and other hydrocarbons.]	266-026-1	65996-91-0	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-046-00-6	Anthracene oil, acid ext.; Anthracene Oil Extract Residue; [A complex combination of hydrocarbons from the base-freed fraction obtained from the distillation of coal tar and boiling in the range of approximately 325 °C to 365 °C (617 °F to 689 °F). It contains predominantly anthracene and phenanthrene and their alkyl derivatives.]	295-274-3	91995-14-1	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-047-00-1	Distillates (coal tar); Heavy Anthracene Oil; [The distillate from coal tar having an approximate distillation range of 100 °C to 450 °C (212 °F to 842 °F). Composed primarily of two to four membered condensed ring aromatic hydrocarbons, phenolic compounds, and aromatic nitrogen bases.]	266-027-7	65996-92-1	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-048-00-7	Distillates (coal tar), pitch, heavy oils; Heavy Anthracene Oil; [The distillate from the distillation of the pitch obtained from bituminous high temperature tar. Composed primarily of tri- and polynuclear aromatic hydrocarbons and boiling in the range of approximately 300 °C to 470 °C (572 °F to 878 °F). The product may also contain heteroatoms.]	295-312-9	91995-51-6	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-049-00-2	Distillates (coal tar), pitch; Heavy Anthracene Oil; [The oil obtained from condensation of the vapors from the heat treatment of pitch. Composed primarily of two- to four-ring aromatic compounds boiling in the range of 200 °C to greater than 400 °C (392 °F to greater than 752 °F).]	309-855-7	101316-49-8	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-050-00-8	Distillates (coal tar), heavy oils, pyrene fraction; Heavy Anthracene Oil Redistillate; [The redistillate obtained from the fractional distillation of pitch distillate boiling in the range of approximately 350 °C to 400 °C (662 °F to 752 °F). Consists predominantly of tri- and polynuclear aromatics and heterocyclic hydrocarbons.]	295-304-5	91995-42-5	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-051-00-3	Distillates (coal tar), pitch, pyrene fraction; Heavy Anthracene Oil Redistillate; [The redistillate obtained from the fractional distillation of pitch distillate and boiling in the range of approximately 380 °C to 410 °C (7160 to 770 °F). Composed primarily of tri- and polynuclear aromatic hydrocarbons and heterocyclic compounds.]	295-313-4	91995-52-7	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-052-00-9	Paraffin waxes (coal), brown-coal high-temp. tar, carbon-treated; Coal Tar Extract; [A complet combination of hydrocarbons obtained by the treatment of lignite carbonization tar with activated carbon for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-296-6	97926-76-6	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-053-00-4	Paraffin waxes (coal), brown-coal high-temp tar, clay-treated; Coal Tar Extract; [A complex combination of hydrocarbons obtained by the treatment of lignite carbonization tar with bentonite for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-297-1	97926-77-7	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-054-00-X	Pitch; Pitch	263-072-4	61789-60-4	Carc. 1B	H350	GHS08 Dgr	H350			H M

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648-055-00-5	Pitch, coal tar, high-temp.; Pitch; [The residue from the distillation of high temperature coal tar. A black solid with an approximate softening point from 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of a complex mixture of three or more membered condensed ring aromatic hydrocarbons.]	266-028-2	65996-93-2	Carc. 1B	H350	GHS08 Dgr	H350			H
648-056-00-0	Pitch, coal tar, high-temp., heat-treated; Pitch; [The heat treated residue from the distillation of high temperature coal tar. A black solid with an approximate softening point from 80 °C to 180 °C (176 °F to 356 °F). Composed primarily of a complex mixture of three or more membered condensed ring aromatic hydrocarbons.]	310-162-7	121575-60-8	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-057-00-6	Pitch, coal tar, high-temp., secondary; Pitch Redistillate; [The residue obtained during the distillation of high boiling fractions from bituminous coal high temperature tar and/or pitch coke oil, with a softening point of 140 °C to 170 °C (284 °F to 392 °F) according to DIN 52025. Composed primarily of tri- and polynuclear aromatic compounds which also contain heteroatoms.]	302-650-3	94114-13-3	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-058-00-1	Residues (coal tar), pitch distn.; Pitch Redistillate; [Residue from the fractional distillation of pitch distillate boiling in the range of approximately 400 °C to 470 °C (752 °F to 846 °F). Composed primarily of polynuclear aromatic hydrocarbons, and heterocyclic compounds.]	295-507-9	92061-94-4	Carc. 1B	H350	GHS08 Dgr	H350			H M

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648-059-00-7	Tar, coal, high-temp., distn. and storage residues; Coal Tar Solids Residue; [Coke- and ash-containing solid residues that separate on distillation and thermal treatment of bituminous coal high temperature tar in distillation installations and storage vessels. Consists predominantly of carbon and contains a small quantity of hetero compounds as well as ash components.]	295-535-1	92062-20-9	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-060-00-2	Tar, coal, storage residues; Coal Tar Solids Residue; [The deposit removed from crude coal tar storages. Composed primarily of coal tar and carbonaceous particulate matter.]	293-764-1	91082-50-7	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-061-00-8	Tar, coal, high-temp., residues; Coal Tar Solids Residue; [Solids formed during the coking of bituminous coal to produce crude bituminous coal high temperature tar. Composed primarily of coke and coal particles, highly aromatized compounds and mineral substances.]	309-726-5	100684-51-3	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-062-00-3	Tar, coal, high-temp., high-solids; Coal Tar Solids Residue; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in the high temperature (greater than 700 °C (1 292 °F)) destructive distillation of coal. Composed primarily of a complex mixture of condensed ring aromatic hydrocarbons with a high solid content of coal-type materials.]	273-615-7	68990-61-4	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-063-00-9	Waste solids, coal-tar pitch coking; Coal Tar Solids Residue; [The combination of wastes formed by the coking of bituminous coal tar pitch. It consists predominantly of carbon.]	295-549-8	92062-34-5	Carc. 1B	H350	GHS08 Dgr	H350			H M

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648-064-00-4	Extract residues (coal), brown; Coal Tar Extract; [The residue from extraction of dried coal.]	294-285-0	91697-23-3	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-065-00-X	Paraffin waxes (coal), brown-coal-high-temp. tar; Coal Tar Extract; [A complex combination of hydrocarbons obtained from lignite carbonization tar by solvent crystallisation (solvent deoiling), by sweating or an adducting process. It consists predominantly of straight and branched chain saturated hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-454-1	92045-71-1	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-066-00-5	Paraffin waxes (coal), brown-coal-high-temp. tar, hydrotreated; Coal Tar Extract; [A complex combination of hydrocarbons obtained from lignite carbonization tar by solvent crystallisation (solvent deoiling), by sweating or an adducting process treated with hydrogen in the presence of a catalyst. It consists predominantly of straight and branched chain saturated hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-455-7	92045-72-2	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-067-00-0	Paraffin waxes (coal), brown-coal high-temp tar, silicic acid-treated; Coal Tar Extract; [A complex combination of hydrocarbons obtained by the treatment of lignite carbonization tar with silicic acid for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-298-7	97926-78-8	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-068-00-6	Tar, coal, low-temp., distn. residues; Tar Oil, intermediate boiling; [Residues from fractional distillation of low temperature coal tar to remove oils that boil in a range up to approximately 300 °C (572 °F). Composed primarily of aromatic compounds.]	309-887-1	101316-85-2	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-069-00-1	Pitch, coal tar, low-temp; Pitch Residue; [A complex black solid or semi-solid obtained from the distillation of a low temperature coal tar. It has a softening point within the approximate range of 40 °C to 180 °C (104 °F to 356 °F). Composed primarily of a complex mixture of hydrocarbons.]	292-651-4	90669-57-1	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-070-00-7	Pitch, coal tar, low-temp., oxidized; Pitch Residue, oxidised; [The product obtained by air-blowing, at elevated temperature, low-temperature coal tar pitch. It has a softening-point within the approximate range of 70 °C to 180 °C (158 °F to 356 °F). Composed primarily of a complex mixture of hydrocarbons.]	292-654-0	90669-59-3	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-071-00-2	Pitch, coal tar, low-temp., heat-treated; Pitch Residue, oxidised; Pitch Residue, heat-treated; [A complex black solid obtained by the heat treatment of low temperature coal tar pitch. It has a softening point within the approximate range of 50 °C to 140 °C (122 °F to 284 °F). Composed primarily of a complex mixture of aromatic compounds.]	292-653-5	90669-58-2	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-072-00-8	Distillates (coal-petroleum), condensed-ring arom; Distillates; [The distillate from a mixture of coal and tar and aromatic petroleum streams having an approximate distillation range of 220 °C to 450 °C (428 °F to 842 °F). Composed primarily of 3- to 4-membered condensed ring aromatic hydrocarbons.]	269-159-3	68188-48-7	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-073-00-3	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polyethylene-polypropylene pyrolysis-derived; Pyrolysis Products; [A complex combination hydrocarbons obtained from mixed coal tar pitch-polyethylene-polypropylene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-956-6	101794-74-5	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-074-00-9	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polyethylene pyrolysis-derived; Pyrolysis Products; [A complex combination of hydrocarbons obtained from mixed coal tar pitch-polyethylene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-957-1	101794-75-6	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-075-00-4	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polystyrene pyrolysis-derived; Pyrolysis Products; [A complex combination of hydrocarbons obtained from mixed coal tar pitch-polystyrene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-958-7	101794-76-7	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-076-00-X	Pitch, coal tar-petroleum; Pitch Residues; [The residue from the distillation of a mixture of coal tar and aromatic petroleum streams. A solid with a softening point from 40 °C to 180 °C (140 °F to 356 °F). Composed primarily of a complex combination of three or more membered condensed ring aromatic hydrocarbons.]	269-109-0	68187-57-5	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-077-00-5	Phenanthrene, distn. residues; Heavy Anthracene Oil Redistillate; [Residue from the distillation of crude phenanthrene boiling in the approximate range of 340 °C to 420 °C (644 °F to 788 °F). It consists predominantly of phenanthrene, anthracene and carbazole.]	310-169-5	122070-78-4	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-078-00-0	Distillates (coal tar), upper, fluorene-free; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained by the crystallization of tar oil. It consists of aromatic polycyclic hydrocarbons, primarily diphenyl, dibenzofuran and acenaphthene.]	284-899-7	84989-10-6	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-079-00-6	Anthracene oil; Anthracene oil; [A complex combination of polycyclic aromatic hydrocarbons obtained from coal tar having an approximate distillation range of 300 °C ot 400 °C (572 °F to 752 °F). Composed primarily of phenanthrene, anthracene and carbazole.]	292-602-7	90640-80-5	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-080-00-1	Residues (coal tar), creosote oil distn.; Wash Oil Redistillate; [The residue from the fractional distillation of wash oil boiling in the approximate range of 270 °C to 330 °C (518 °F to 626 °F). It consists predominantly of dinuclear aromatic and heterocyclic hydrocarbons.]	295-506-3	92061-93-3	Carc. 1B	H350	GHS08 Dgr	H350			H
648-081-00-7	Tar, coal; Coal tar; [The by-product from the destructive distillation of coal. Almost black semisolid. A complex combination of aromatic hydrocarbons, phenolic compounds, nitrogen bases and thiophene.]	232-361-7	8007-45-2	Carc. 1A	H350	GHS08 Dgr	H350			H
648-082-00-2	Tar, coal, high-temp.; Coal tar; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in the high temperature (greater than 700 °C (1 292 °F)) destructive distillation of coal. A black viscous liquid denser than water. Composed primarily of a complex mixture of condensed ring aromatic hydrocarbons. May contain minor amounts of phenolic compounds and aromatic nitrogen bases.]	266-024-0	65996-89-6	Carc. 1A	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-083-00-8	Tar, coal, low-temp.; Coal oil; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in low temperature (less than 700 °C (1292 °F)) destructive distillation of coal. A black viscous liquid denser than water. Composed primarily of condensed ring aromatic hydrocarbons, phenolic compounds, aromatic nitrogen bases, and their alkyl derivatives.]	266-025-6	65996-90-9	Carc. 1A	H350	GHS08 Dgr	H350			H
648-084-00-3	Distillates (coal), coke-oven light oil, naphthalene cut; Naphthalene Oil; [The complex combination of hydrocarbons obtained from prefractionation (continuous distillation) of coke oven light oil. It consists predominantly of naphthalene, coumarone and indene and boils above 148 °C (298 °F).]	285-076-5	85029-51-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-085-00-9	Distillates (coal tar), naphthalene oils; Naphthalene Oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists primarily of aromatic and other hydrocarbons, phenolic compounds and aromatic nitrogen compounds and distills in the approximate range of 200 °C to 250 °C (392 °F to 482 °F).]	283-484-8	84650-04-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-086-00-4	Distillates (coal tar), naphthalene oils, naphthalene-low; Naphthalene Oil Redistillate; [A complex combination of hydrocarbons obtained by crystallization of naphthalene oil. Composed primarily of naphthalene, alkyl naphthalenes and phenolic compounds.]	284-898-1	84989-09-3	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-087-00-X	Distillates (coal tar), naphthalene oil crystn. mother liquor; Naphthalene Oil Redistillate; [A complex combination of organic compounds obtained as a filtrate from the crystallization of the naphthalene fraction from coal tar and boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). Contains chiefly naphthalene, thionaphthene and alkylnaphthalenes.]	295-310-8	91995-49-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-088-00-5	Extract residues (coal), naphthalene oil, alk.; Naphthalene Oil Extract Residue; [A complex combination of hydrocarbons obtained from the alkali washing of naphthalene oil to remove phenolic compounds (tar acids). It is composed of naphthalene and alkyl naphthalenes.]	310-166-9	121620-47-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-089-00-0	Extract residues (coal), naphthalene oil, alk., naphthalene-low; Naphthalene Oil Extract Residue; [A complex combination of hydrocarbons remaining after the removal of naphthalene from alkali-washed naphthalene oil by a crystallization process. It is composed primarily of naphthalene and alkyl naphthalenes.]	310-167-4	121620-48-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-090-00-6	Distillates (coal tar), naphthalene oils, naphthalene-free, alk. exts.; Naphthalene Oil Extract Residue; [The oil remaining after the removal of phenolic compounds (tar acids) from drained naphthalene oil by an alkali wash. Composed primarily of naphthalene and alkyl naphthalenes.]	292-612-1	90640-90-7	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-091-00-1	Extract residues (coal), naphthalene oil alk., distn. overheads; Naphthalene Oil Extract Residue; [The distillation from alkali-washed naphthalene oil having an approximate distillation range of 180 °C to 220 °C (356 °F to 428 °F). Composed primarily of naphthalene, alkylbenzenes, indene and indan.]	292-627-3	90641-04-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-092-00-7	Distillates (coal tar), naphthalene oils, methylnaphthalene fraction; Methylnaphthalene Oil; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of substituted two ring aromatic hydrocarbons and aromatic nitrogen bases boiling in the range of approximately 225 °C to 255 °C (437 °F to 491 °F).]	309-985-4	101896-27-9	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-093-00-2	Distillates (coal tar), naphthalene oils, indole-methylnaphthalene fraction; Methylnaphthalene Oil; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of indole and methylnaphthalene boiling in the range of approximately 235 °C to 255 °C (455 °F to 491 °F).]	309-972-3	101794-91-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-094-00-8	Distillates (coal tar), naphthalene oils, acid exts.; Methylnaphthalene Oil Extract Residue; [A complex combination of hydrocarbons obtained by debasing the methylnaphthalene fraction obtained by the distillation of coal tar and boiling in the range of approximately 230 °C to 255 °C (446 °F to 491 °F). Contains chiefly 1(2)-methylnaphthalene, naphthalene, dimethylnaphthalene and biphenyl.]	295-309-2	91995-48-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-095-00-3	Extract residues (coal), naphthalene oil alk., distn. residues; Methylnaphthalene Oil Extract Residue; [The residue from the distillation of alkali-washed naphthalene oil having an approximate distillation range of 220 °C to 300 °C (428 °F to 572 °F). Composed primarily of naphthalene, alkylnaphthalenes and aromatic nitrogen bases.]	292-628-9	90641-05-7	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-096-00-9	Extract oils (coal), acidic, tar-base free; Methylnaphthalene Oil Extract Residue; [The extract oil boiling in the range of approximately 220 °C to 265 °C (428 °F to 509 °F) from coal tar alkaline extract residue produced by an acidic wash such as aqueous sulfuric acid after distillation to remove tar bases. Composed primarily of alkylnaphthalenes.]	284-901-6	84989-12-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-097-00-4	Distillates (coal tar), benzole fraction, distn. residues; Wash Oil; [A complex combination of hydrocarbons obtained from the distillation of crude benzole (high temperature coal tar). It may be a liquid with the approximate distillation range of 150 °C to 300 °C (302 °F to 572 °F) or a semi-solid or solid with a melting point up to 70 °C (158 °F). It is composed primarily of naphthalene and alkyl naphthalenes.]	310-165-3	121620-46-0	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-098-00-X	Creosote oil, acenaphthene fraction; Wash Oil; [A complex combination of hydrocarbons produced by the distillation of coal tar and boiling in the range of approximately 240 °C to 280 °C (464 °F to 536 °F). Composed primarily of acenaphthene, naphthalene and alkyl naphthalene.]	292-605-3	90640-84-9	Carc. 1B	H350	GHS08 Dgr	H350			H

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648-099-00-5	Creosote oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists primarily of aromatic hydrocarbons and may contain appreciable quantities of tar acids and tar bases. It distills at the approximate range of 200 °C to 325 °C (392 °F to 617 °F).]	263-047-8	61789-28-4	Carc. 1B	H350	GHS08 Dgr	H350			H
648-100-00-9	Creosote oil, high-boiling distillate; Wash Oil; [The high-boiling distillation fraction obtained from the high temperature carbonization of bituminous coal which is further refined to remove excess crystalline salts. It consists primarily of creosote oil with some of the normal polynuclear aromatic salts, which are components of coal tar distillates, removed. It is crystal free at approximately 5 °C (41 °F).]	274-565-9	70321-79-8	Carc. 1B	H350	GHS08 Dgr	H350			H
648-101-00-4	Creosote; [The distillate of coal tar produced by the high temperature carbonization of bituminous coal. It consists primarily of aromatic hydrocarbons, tar acids and tar bases.]	232-287-5	8001-58-9	Carc. 1B	H350	GHS08 Dgr	H350			H
648-102-00-X	Extract residues (coal), creosote oil acid; Wash Oil Extract Residue; [A complex combination of hydrocarbons from the base-freed fraction from the distillation of coal tar, boiling in the range of approximately 250 °C to 280 °C (482 °F to 536 °F). It consists predominantly of biphenyl and isomeric diphenylnaphthalenes.]	310-189-4	122384-77-4	Carc. 1B	H350	GHS08 Dgr	H350			H
648-103-00-5	Anthracene oil, anthracene paste; Anthracene Oil Fraction; [The anthracene-rich solid obtained by the crystallization and centrifuging of anthracene oil. It is composed primarily of anthracene, carbazole and phenanthrene.]	292-603-2	90640-81-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-104-00-0	Anthracene oil, anthracene-low; Anthracene Oil Fraction; [The oil remaining after the removal, by a crystallization process, of an anthracene-rich solid (anthracene paste) from anthracene oil. It is composed primarily of two, three and four membered aromatic compounds.]	292-604-8	90640-82-7	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-105-00-6	Residues (coal tar), anthracene oil distn.; Anthracene Oil Fraction; [The residue from the fraction distillation of crude anthracene boiling in the approximate range of 340 °C to 400 °C (644 °F to 752 °F). It consists predominantly of tri- and polynuclear aromatic and heterocyclic hydrocarbons.]	295-505-8	92061-92-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-106-00-1	Anthracene oil, anthracene paste, anthracene fraction; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by the crystallization of anthracene oil from bituminous high temperature tar and boiling in the range of 330 °C to 350 °C (626 °F to 662 °F). It contains chiefly anthracene, carbazole and phenanthrene.]	295-275-9	91995-15-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-107-00-7	Anthracene oil, anthracene paste, carbazole fraction; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by crystallization of anthrancene oil from bituminous coal high temperature tar and boiling in the approximate range of 350 °C to 360 °C (662 °F to 680 °F). It contains chiefly anthacene, carbazole and phenanthrene.]	295-276-4	91995-16-3	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-108-00-2	Anthracene oil, anthracene paste, distn. lights; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by crystallization of anthracene oil from bituminous light temperature tar and boiling in the range of approximately 290 °C to 340 °C (554 °F to 644 °F). It contains chiefly trinuclear aromatics and their dihydro derivatives.]	295-278-5	91995-17-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-109-00-8	Tar oils, coal, low-temp.; Tar Oil, high boiling; [A distillate from low-temperature coal tar. Composed primarily of hydrocarbons, phenolic compounds and aromatic nitrogen bases boiling in the range of approximately 160 °C to 340 °C (320 °F to 644 °F).]	309-889-2	101316-87-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-110-00-3	Extract residues (coal), low temp. coal atar alk.; [The residue from low temperature coal tar oils after an alkaline wash, such as aqueous sodium hydroxide, to remove crude coal tar acids. Composed primarily of hydrocarbons and aromatic nitrogen bases.]	310-191-5	122384-78-5	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-111-00-9	Phenols, ammonia liquor ext.; Alkaline Extract; [The combination of phenols extracted, using isobutyl acetate, from the ammonia liquor condensed from the gas evolved in low-temperature (less than 700 °C (1292 °F)) destructive distillation of coal. It consists predominantly of a mixture of monohydric and dihydric phenols.]	284-881-9	84988-93-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-112-00-4	Distillates (coal tar), light oils, alk. exts.; Alkaline Extract; [The aqueous extract from carbolic oil produced by an alkaline wash such as aqueous sodium hydroxide. Composed primarily of the alkali salts of various phenolic compounds.]	292-610-0	90640-88-3	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-113-00-X	Extracts, coal tar oil alk.; Alkaline Extract; [The extract from coal tar oil produced by an alkaline wash such as aqueous sodium hydroxide. Composed primarily of the alkali salts of various phenolic compounds.]	266-017-2	65996-83-0	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-114-00-5	Distillates (coal tar), naphthalene oils, alk. exts.; Alkaline Extract; [The aqueous extract from naphthalene oil produced by an alkaline wash such as aqueous sodium hydroxid. Composed primarily of the alkali salts of various phenolic compounds.]	292-611-6	90640-89-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-115-00-0	Extract residues (coal), tar oil alk., carbonated, limed; Crude Phenols; [The product obtained by treatment of coal tar oil alkaline extract with CO_2 and CaO. Composed primarily of $CaCO_3$, $Ca(OH)_2$, Na_2CO_3 and other organic and inorganic impurities.]	292-629-4	90641-06-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-116-00-6	Tar acids, coal, crude; Crude Phenols; [The reaction product obtained by neutralizing coal tar oil alkaline extract with an acidic solution, such as aqueous sulfuric acid, or gaseous carbon dioxide, to obtain the free acids. Composed primarily of tar acids such as phenol, cresols, and xylenols.]	266-019-3	65996-85-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-117-00-1	Tar acids, brown-coal, crude; Crude Phenols; [An acidified alkaline extract of brown coal tar distillate. Composed primarily of phenol and phenol homologs.]	309-888-7	101316-86-3	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-118-00-7	Tar acids, brown-coal gasification; Crude Phenols; [A complex combination of organic compounds obtained from brown coal gasification. Composed primarily of C_{6-10} hydroxy aromatic phenols and their homologs.]	295-536-7	92062-22-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-119-00-2	Tar acids, distn. residues; Distillate Phenols; [A residue from the distillation of crude phenol from coal. It consists predominantly of phenols having carbon numbers in the range of C_8 through C_{10} with a softening point of 60 °C to 80 °C (140 °F to 176 °F).]	306-251-5	96690-55-0	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-120-00-8	Tar acids, methylphenol fraction; Distillate Phenols; [The fraction of tar acid rich in 3- and 4-methylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-892-9	84989-04-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-121-00-3	Tar acids, polyalkylphenol fraction; Distillate Phenols; [The fraction of tar aicds, recovered by distillation of low-temperature coal tar crude tar acids, having an approximate boiling range of 225 °C to 320 °C (437 °F to 608 °F). Composed primarily of polyalkylphenols.]	284-893-4	84989-05-9	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-122-00-9	Tar acids, xylenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 2,4- and 2,5-dimethylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-895-5	84989-06-0	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-123-00-4	Tar acids, ethylphenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 3- and 4-ethylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-891-3	84989-03-7	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-124-00-X	Tar acids, 3,5-xylenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 3,5-dimethylphenol, recovered by distillation of low-temperature coal tar acids.]	284-896-0	84989-07-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-125-00-5	Tar acids, residues, distillates, first-cut; Distillate Phenols; [The residue from the distillation in the range of 235 °C to 355 °C (481 °F to 697 °F) of light carbolic oil.]	270-713-1	68477-23-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-126-00-0	Tar acids, cresylic, residues; Distillate Phenols; [The residue from crude coal tar acids after removal of phenol, cresols, xylenols and any higher boiling phenols. A black solid with a melting point approximately 80 °C (176 °F). Composed primarily of polyalkyphenols, resin gums, and inorganic salts.]	271-418-0	68555-24-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-127-00-6	Phenols, C_{9-11}; Distillate Phenols	293-435-2	91079-47-9	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-128-00-1	Tar acids, cresylic; Distillate Phenols; [A complex combination of organic compounds obtained from brown coal and boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). It contains chiefly phenols and pyridine bases.]	295-540-9	92062-26-5	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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648-129-00-7	Tar acids, brown-coal, C_2-alkylphenol fraction; Distillate Phenols; [The distillate from the acidification of alkaline washed lignite tar distillate boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). Composed primarily of *m*- and *p*-ethylphenol as well as cresols and xylenols.]	302-662-9	94114-29-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-130-00-2	Extract oils (coal), naphthalene oils; Acid Extract; [The aqueous extract produced by an acidic wash of alkali-washed napthalene oil. Composed primarily of acid salts of various aromatic nitrogen bases including pyridine, quinoline and their alkyl derivatives.]	292-623-1	90641-00-2	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-131-00-8	Tar bases, quinoline derivs.; Distillate Bases	271-020-7	68513-87-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-132-00-3	Tar bases, coal, quinoline derivs. fraction; Distillate Bases	274-560-1	70321-67-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-133-00-9	Tar bases, coal, distn. residues; Distillate Bases; [The distillation residue remaining after the distillation of the neutralized, acid-extracted base-containing tar fractions obtained by the distillation of coal tars. It contains chiefly aniline, collidines, quinoline and quinoline derivatives and toluidines.]	295-544-0	92062-29-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-134-00-4	Hydrocarbon oils, arom., mixed with polyethylene and polypropylene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of a polyethylene/polypropylene mixture with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of approximately 70 °C to 120 °C (158 °F to 248 °F).]	309-745-9	100801-63-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-135-00-X	Hydrocarbon oils, arom., mixed with polyethylene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of polyethylene with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of 70 °C to 120 °C (158 °F to 248 °F).]	309-748-5	100801-65-8	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-136-00-5	Hydrocarbon oils, arom., mixed with polystyrene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of polystyrene with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of approximately 70 °C to 210 °C (158 °F to 410 °F).]	309-749-0	100801-66-9	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-137-00-0	Extract residues (coal), tar oil alk., naphthalene distn. residues; Naphthalene Oil Extract Residue; [The residue obtained from chemical oil extracted after the removal of naphthalene by distillation composed primarily of two to four membered condensed ring aromatic hydrocarbons and aromatic nitrogen bases.]	277-567-8	73665-18-6	Carc. 1B	H350	GHS08 Dgr	H350			HJM

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-138-00-6	Creosote oil, low-boiling distillate; Wash Oil; [The low-boiling distillation fraction obtained from the high temperature carbonization of bituminous coal, which is further refined to remove excess crystalline salts. It consists primarily of creosote oil with some of the normal polynuclear aromatic salts, which are components of coal tar distillate, removed. It is crystal free at approximately 38 °C (100 °F).]	274-566-4	70321-80-1	Carc. 1B	H350	GHS08 Dgr	H350			H
648-139-00-1	Tar acids, cresylic, sodium salts, caustic solns.; Alkaline Extract	272-361-4	68815-21-4	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-140-00-7	Extract oils (coal), tar base; Acid Extract; [The extract from coal tar oil alkaline extract residue produced by an acidic wash such as aqueous sulfuric acid after distillatin to remove naphthalene. Composed primarily of the acid salts of various aromatic nitrogen bases including pyridine, quinoline, and their alkyl derivatives.]	266-020-9	65996-86-3	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-141-00-2	Tar bases, coal, crude; Crude Tar Bases; [The reaction product obtained by neutralizing coal tar base extract oil with an alkaline solution, such as aqueous sodium hydroxide, to obtain the free bases. Composed primarily of such organic bases as acridine, phenanthridine, pyridine, quinoline and their alkyl derivatives.]	266-018-8	65996-84-1	Carc. 1B	H350	GHS08 Dgr	H350			HJM
648-142-00-8	Residues (coal), liq. solvent extn.; [A cohesive powder composed of coal mineral matter and undissolved coal remaining after extraction of coal by a liquid solvent.]	302-681-2	94114-46-2	Carc. 1B	H350	GHS08 Dgr	H350			H M

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-143-00-3	Coal liquids, liq. solvent extn. soln.; [The product obtained by filtration of coal mineral matter and undissolved coal from coal extract solution produced by digesting coal in a liquid solvent. A black, viscous, highly complex liquid combination composed primarily of aromatic and partly hydro-genated aromatic hydrocarbons, aromatic nitrogen compounds, aromatic sulfur compounds, phenolic and other aromatic oxygen compounds and their alkyl derivatives.]	302-682-8	94114-47-3	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-144-00-9	Coal liquids, liq. solvent extn.; [The substantially solvent-free product obtained by the distillation of the solvent from filtered coal extract solution produced by digesting coal in a liquid solvent. A black semi-solid, composed primarily of a complex combination of condensed-ring aromatic hydrocarbons, aromatic nitrogen compounds, aromatic sulfur compounds, phenolic compounds and other aromatic oxygen compounds, and their alkyl derivatives.]	302-683-3	94114-48-4	Carc. 1B	H350	GHS08 Dgr	H350			H M
648-145-00-4	Tar brown-coal; [An oil distilled from brown-coal tar. Composed primarily of aliphatic, naphthenic and one- to three-ring aromatic hydrocarbons, their alkyl derivates, heteroaromatics and one- and two-ring phenols boiling in the range of approximately 150 °C to 360 °C (302 °F to 680 °F).]	309-885-0	101316-83-0	Carc. 1A	H350	GHS08 Dgr	H350			H
648-146-00-X	Tar, brown-coal, low-temp.; [A tar obtained from low temperature carbonization and low temperature gasification of brown coal. Composed primarily of aliphatic, naphthenic and cyclic aromatic hydrocarbons, heteroaromatic hydrocarbons and cyclic phenols.]	309-886-6	101316-84-1	Carc. 1A	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-147-00-5	Light oil (coal), coke-oven; Crude benzole; [The volatile organic liquid extracted from the gas evolved in the high temperature (greater than 700 °C (1292 °F)) destructive distillation of coal. Composed primarily of benzene, toluene, and xylenes. May contain other minor hydrocarbon constituents.]	266-012-5	65996-78-3	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-148-00-0	Distillates (coal), liq. solvent extn., primary; [The liquid product of condensation of vapors emitted during the digestion of coal in a liquid solvent and boiling in the range of approximately 30 °C to 300 °C (86 °F to 572 °F). Composed primarily of partly hydrogenated condensed-ring aromatic hydrocarbons, aromatic compounds containing nitrogen, oxygen and sulfur, and their alkyl derivatives having carbon numbers predominantly in the range of C_4 through C_{14}.]	302-688-0	94114-52-0	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-149-00-6	Distillates (coal), solvent extn., hydrocracked; [Distillate obtained by hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction process and boiling in the range of approximately 30 °C to 300 °C (86 °F to 572 °F). Composed primarily of aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes with carbon numbers predominantly in the range of C_4 through C_{14}. Nitrogen, sulfur and oxygen-containing aromatic and hydrogenated aromatic compounds are also present.]	302-689-6	94114-53-1	Carc. 1B	H350	GHS08 Dgr	H350			H J

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648-150-00-1	Naphtha (coal), solvent extn., hydrocracked; [Fraction of the distillate obtained by hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes with carbon numbers predominantly in the range of C_4 to C_9. Nitrogen, sulfur and oxygen-containing aromatic and hydrogenated aromatic compounds are also present.]	302-690-1	94114-54-2	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-151-00-7	Gasoline, coal solvent extn., hydrocracked naphtha; [Motor fuel produced by the reforming of the refined naphtha fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of aromatic and naphthenic hydrocarbons, their alkyl derivatives and alkyl hydrocarbons having carbon numbers in the range of C_4 through C_9.]	302-691-7	94114-55-3	Carc. 1B	H350	GHS08 Dgr	H350			H
648-152-00-2	Distillates (coal), solvent extn., hydrocracked middle; [Distillate obtained from the hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 300 °C (356 °F to 572 °F). Composed primarily of two-ring aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes having carbon numbers predominantly in the range of C_9 through C_{14}. Nitrogen, sulfur and oxygen-containing compounds are also present.]	302-692-2	94114-56-4	Carc. 1B	H350	GHS08 Dgr	H350			H J

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-153-00-8	Distillates (coal), solvent extn., hydrocracked hydrogenated middle; [Distillate from the hydrogenation of hydrocracked middle distillate from coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 280 °C (356 °F to 536 °F). Composed primarily of hydrogenated two- ring carbon compounds and their alkyl derivatives having carbon numbers predominantly in the range of C_9 through C_{14}.]	302-693-8	94114-57-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
648-154-00-3	Fuels, jet aircraft, coal solvent extn., hydrocracked hydrogenated; [Jet engine fuel produced by hydrogenation of the middle distillate fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 225 °C (356 °F to 473 °F). Composed primarily of hydrogenated two-ring hydrocarbons and their alkyl derivatives having carbon numbers predominantly in the range of C_{10} through C_{12}.]	302-694-3	94114-58-6	Carc. 2	H351	GHS08 Wng	H350			H
648-155-00-9	Fuels, diesel, coal solvent extn., hydrocracked hydrogenated; [Diesel engine fuel produced by the hydrogenation of the middle distillate fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 200 °C to 280 °C (392 °F to 536 °F). Composed primarily of hydrogenated two-ring hydrocarbons and their alkyl derivatives having carbon numbers predominantly in the range of C_{11} through C_{14}.]	302-695-9	94114-59-7	Carc. 2	H351	GHS08 Wng	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
648-156-00-4	Light oil (coal), semi-coking process; Fresh oil; [The volatile organic liquid condensed from the gas evolved in the low temperature (less than 700 °C (1292 °F) destructive distillation of coal. Composed primarily of C_{6-10} hydrocarbons.]	292-635-7	90641-11-5	Carc. 1B	H350	GHS08 Dgr	H350			H J
649-001-00-3	Extracts (petroleum), light naphthenic distillate solvent	265-102-1	64742-03-6	Carc. 1B	H350	GHS08 Dgr	H350			H
649-002-00-9	Extracts (petroleum), heavy paraffinic distillate solvent	265-103-7	64742-04-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-003-00-4	Extracts (petroleum), light paraffinic distillate solvent	265-104-2	64742-05-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-004-00-X	Extracts (petroleum), heavy naphthenic distillate solvent	265-111-0	64742-11-6	Carc. 1B	H350	GHS08 Dgr	H350			H
649-005-00-5	Extracts (petroleum), light vacuum gas oil solvent	295-341-7	91995-78-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-006-00-0	hydrocarbons C_{26-55}, arom-rich	307-753-7	97722-04-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-007-00-6	fatty acids, tall-oil, reaction products with iminodiethanol and boric acid	400-160-5	—	Skin Irrit. 2 Aquatic Chronic 2	H315 H411	GHS07 GHS09 Wng	H315 H411			
649-008-00-1	Residues (petroleum), atm. tower; Heavy Fuel oil; [A complex residuum from the atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-045-2	64741-45-3	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-009-00-7	Gas oils (petroleum), heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and boiling in the range of approximately 350 °C to 600 °C (662 °F to 1112 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	265-058-3	64741-57-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-010-00-2	Distillates (petroleum), heavy catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{35} and boiling in the range of approximately 260 °C to 500 °C (500 °F to 932 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-063-0	64741-61-3	Carc. 1B	H350	GHS08 Dgr	H350			H
649-011-00-8	Clarified oils (petroleum), catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-064-6	64741-62-4	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-012-00-3	Residues (petroleum), hydrocracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the products of a hydrocracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F).]	265-076-1	64741-75-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-013-00-9	Residues (petroleum), thermal cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the product from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-081-9	64741-80-6	Carc. 1B	H350	GHS08 Dgr	H350			H
649-014-00-4	Distillates (petroleum), heavy thermal cracked; Heavy Fuel oil; [A complex combination of hydrocarbons from the distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{36} and boiling in the range of approximately 260 °C to 480 °C (500 °F to 896 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-082-4	64741-81-7	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-015-00-X	Gas oils (petroleum), hydrotreated vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{50} and boiling in the range of approximately 230 °C to 600 °C (446 °F to 1112 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-162-9	64742-59-2	Carc. 1B	H350	GHS08 Dgr	H350			H
649-016-00-5	Residues (petroleum), hydrodesulfurized atmospheric tower; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating an atmospheric tower residuum with hydrogen in the presence of a catalyst under conditions primarily to remove organic sulfur compounds. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-181-2	64742-78-5	Carc. 1B	H350	GHS08 Dgr	H350			H
649-017-00-0	Gas oils (petroleum), hydrodesulfurized heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and boiling in the range of approximately 350 °C to 600 °C (662 °F to 1112 °C). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-189-6	64742-86-5	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-018-00-6	Residues (petroleum), steam-cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained as the residual fraction from the distillation of the products of a steam cracking process (including steam cracking to produce ethylene). It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly greater than C_{14} and boiling above approximately 260 °C (500 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-193-8	64742-90-1	Carc. 1B	H350	GHS08 Dgr	H350			H
649-019-00-1	Residues (petroleum), atmospheric; Heavy Fuel oil; [A complex residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{11} and boiling above approximately 200 °C (392 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	269-777-3	68333-22-2	Carc. 1B	H350	GHS08 Dgr	H350			H
649-020-00-7	Clarified oils (petroleum), hydrodesulfurized catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating catalytic cracked clarified oil with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	269-782-0	68333-26-6	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-021-00-2	Distillates (petroleum), hydrodesulfurized intermediate catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating intermediate catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{30} and boiling in the range of approximately 205 °C to 450 °C (401 °F to 842 °F). It contains a relatively large proportion of tricyclic aromatic hydrocarbons.]	269-783-6	68333-27-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-022-00-8	Distillates (petroleum), hydrodesulfurized heavy catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treatment of heavy catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{35} and boiling in the range of approximately 260 °C to 500 °C (500 °F to 932 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	269-784-1	68333-28-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-023-00-3	Fuel oil, residues-straight-run gas oils, high-sulfur; Heavy Fuel oil	270-674-0	68476-32-4	Carc. 1B	H350	GHS08 Dgr	H350			H
649-024-00-9	Fuel oil, residual; Heavy Fuel oil; [The liquid product from various refinery streams, usually residues. The composition is complex and varies with the source of the crude oil.]	270-675-6	68476-33-5	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-025-00-4	Residues (petroleum), catalytic reformer fractionator residue distn.; Heavy Fuel oil; [A complex residuum from the distillation of catalytic reformer fractionator residue. It boils approximately above 399 °C (750 °F).]	270-792-2	68478-13-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-026-00-X	Residues (petroleum), heavy coker gas oil and vacuum gas oil; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of heavy coker gas oil and vacuum gas oil. It predominantly consists of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-796-4	68478-17-1	Carc. 1B	H350	GHS08 Dgr	H350			H
649-027-00-5	Residues (petroleum), heavy coker and light vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of heavy coker gas oil and light vacuum gas oil. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-983-0	68512-61-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-028-00-0	Residues (petroleum), light vacuum; Heavy Fuel oil; [A complex residuum from the vacuum distillation of the residuum from the atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-984-6	68512-62-9	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-029-00-6	Residues (petroleum), steam-cracked light; Heavy Fuel oil; [A complex residuum from the distillation of the products from a steam-cracking process. It consists predominantly of aromatic and unsaturated hydrocarbons having carbon numbers greater than C_7 and boiling in the range of approximately 101 °C to 555 °C (214 °F to 1030 °F).]	271-013-9	68513-69-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-030-00-1	Fuel oil, No 6; Heavy Fuel oil; [A distillate oil having a minimum viscosity of 900 SUS at 37.7 °C (100 °F) to a maximum of 9000 SUS at 37.7 °C (100 °F).]	271-384-7	68553-00-4	Carc. 1B	H350	GHS08 Dgr	H350			H
649-031-00-7	Residues (petroleum), topping plant, low-sulfur; Heavy Fuel oil; [A low-sulfur complex combination of hydrocarbons produced as the residual fraction from the topping plant distillation of crude oil. It is the residuum after the straight-run gasoline cut, kerosene cut and gas oil cut have been removed.]	271-763-7	68607-30-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-032-00-2	Gas oils (petroleum), heavy atmospheric; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{35} and boiling in the range of approximately 121 °C to 510 °C (250 °F to 950 °F).]	272-184-2	68783-08-4	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-033-00-8	Residues (petroleum), coker scrubber, Condensed-ring-arom.-contg.; Heavy Fuel oil; [A very complex combination of hydrocarbons produced as the residual fraction from the distillation of vaccum residuum and the products from a thermal cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed rind aromatic hydrocarbons.]	272-187-9	68783-13-1	Carc. 1B	H350	GHS08 Dgr	H350			H
649-034-00-3	Distillates (petroleum), petroleum residues vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from the atmospheric distillation of crude oil.]	273-263-4	68955-27-1	Carc. 1B	H350	GHS08 Dgr	H350			H
649-035-00-9	Residues (petroleum), steam-cracked, resinous; Heavy Fuel oil; [A complex residuum from the distillation of steam-cracked petroleum residues.]	273-272-3	68955-36-2	Carc. 1B	H350	GHS08 Dgr	H350			H
649-036-00-4	Distillates (petroleum), intermediate vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum, distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{42} and boiling in the range of approximately 250 °C to 545 °C (482 °F to 1013 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	274-683-0	70592-76-6	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-037-00-X	Distillates (petroleum), light vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{35} and boiling in the range of approximately 250 °C to 545 °C (482 °F to 1013 °F).]	274-684-6	70592-77-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-038-00-5	Distillates (petroleum), vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having numbers predominantly in the range of C_{15} through C_{50} and boiling in the range of approximately 270 °C to 600 °C (518 °F to 1112 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	274-685-1	70592-78-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-039-00-0	Gas oils (petroleum), hydrodesulfurized coker heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by hydrodesulfurization of heavy coker distillate stocks, It consists predominantly of hydrocarbons having carbon numbers predominantly in the range C_{18} to C_{44} and boiling in the range of approximately 304 °C to 548 °C (579 °F to 1018 °F). Likely to contain 5 % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	285-555-9	85117-03-9	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-040-00-6	Residues (petroleum), steam-cracked, distillates; Heavy Fuel oil; [A complex combination of hydrocarbons obtained during the production of refined petroleum tar by the distillation of steam cracked tar. It consists predominantly of aromatic and other hydrocarbons and organic sulfur compounds.]	292-657-7	90669-75-3	Carc. 1B	H350	GHS08 Dgr	H350			H
649-041-00-1	Residues (petroleum), vacuum, light; Heavy Fuel oil; [A complex residuum from the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{24} and boiling above approximately 390 °C (734 °F).]	292-658-2	90669-76-4	Carc. 1B	H350	GHS08 Dgr	H350			H
649-042-00-7	Fuel oil, heavy, high-sulfur; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the distillation of crude petroleum. It consists predominantly of aliphatic, aromatic and cycloaliphatic hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	295-396-7	92045-14-2	Carc. 1B	H350	GHS08 Dgr	H350			H
649-043-00-2	Residues (petroleum), catalytic cracking; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of the products from a catalytic cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{11} and boiling above approximately 200 °C (392 °F).]	295-511-0	92061-97-7	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-044-00-8	Distillates (petroleum), intermediate catalytic cracked, thermally degraded; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process which has been used as a heat transfer fluid. It consists predominantly of hydrocarbons boiling in the range of approximately 220 °C to 450 °C (428 °F to 842 °F). This stream is likely to contain organic sulfur compounds.]	295-990-6	92201-59-7	Carc. 1B	H350	GHS08 Dgr	H350			H
649-045-00-3	Residual oils (petroleum); Heavy Fuel oil; [A complex combination of hydrocarbons, sulfur compounds and metal-containing organic compounds obtained as the residue from refinery fractionation cracking processes. It produces a finished oil with a viscosity above 2cSt. at 100 °C.]	298-754-0	93821-66-0	Carc. 1B	H350	GHS08 Dgr	H350			H
649-046-00-9	Residues, steam cracked, thermally treated; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the treatment and distillation of raw steam-cracked naphtha. It consists predominantly of unsaturated hydrocarbons boiling in the range above approximately 180 °C (356 °F).]	308-733-0	98219-64-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-047-00-4	Distillates (petroleum), hydrodesulfurized full-range middle; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating a petroleum stock with hydrogen. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F).]	309-863-0	101316-57-8	Carc. 1B	H350	GHS08 Dgr	H350			H

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649-048-00-X	Residues (petroleum), catalytic reformer fractionator; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the product from a catalytic reforming process. It consists of predominantly aromatic hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{25} and boiling in the range of approximately 160 °C to 400 °C (320 °F to 725 °F). This stream is likely to contain 5 wt. % or more of 4- or 6-membered condensed ring aromatic hydrocarbons.]	265-069-3	64741-67-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-049-00-5	Petroleum; Crude oil; [A complex combination of hydrocarbons, It consists predominantly of aliphatic, alicyclic and aromatic hydrocarbons. It may also contain small amounts of nitrogen, oxygen and sulfur compounds. This category encompasses light, medium, and heavy petroleums, as well as the oils extended from tar sands. Hydrocarbonaceous materials requiring major chemical changes for their recovery or conversion to petroleum refinery feedstocks such as crude shale oils; upgraded shale oils and liquid coal fuels are not included in this definition.]	232-298-5	8002-05-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-050-00-0	Distillates (petroleum), light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated aliphatic hydrocarbons normally present in this distillation range of crude oil.]	265-051-5	64741-50-0	Carc. 1A	H350	GHS08 Dgr	H350			H

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649-051-00-6	Distillates (petroleum), heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated aliphatic hydrocarbons.]	265-052-0	64741-51-1	Carc. 1A	H350	GHS08 Dgr	H350			H
649-052-00-1	Distillates (petroleum), light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-053-6	64741-52-2	Carc. 1A	H350	GHS08 Dgr	H350			H
649-053-00-7	Distillates (petroleum), heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-054-1	64741-53-3	Carc. 1A	H350	GHS08 Dgr	H350			H

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649-054-00-2	Distillates (petroleum), acid-treated heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-117-3	64742-18-3	Carc. 1A	H350	GHS08 Dgr	H350			H
649-055-00-8	Distillates (petroleum), acid-treated light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-118-9	64742-19-4	Carc. 1A	H350	GHS08 Dgr	H350			H
649-056-00-3	Distillates (petroleum), acid-treated heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of a least 100 SUS at 100 °F (19cSt at 40 °C).]	265-119-4	64742-20-7	Carc. 1A	H350	GHS08 Dgr	H350			H

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649-057-00-9	Distillates (petroleum), acid-treated light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-121-5	64742-21-8	Carc. 1A	H350	GHS08 Dgr	H350			H
649-058-00-4	Distillates (petroleum), chemically neutralized heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained from a treating process to remove acidic materials. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of aliphatic hydrocarbons.]	265-127-8	64742-27-4	Carc. 1A	H350	GHS08 Dgr	H350			H
649-059-00-X	Distillates (petroleum), chemically neutralized light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-128-3	64742-28-5	Carc. 1A	H350	GHS08 Dgr	H350			H

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649-060-00-5	Distillates (petroleum), chemically neutralized heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-135-1	64742-34-3	Carc. 1A	H350	GHS08 Dgr	H350			H
649-061-00-0	Distillates (petroleum), chemically neutralized light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS a 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-136-7	64742-35-4	Carc. 1A	H350	GHS08 Dgr	H350			H
649-062-00-6	Gases (petroleum), catalytic cracked naphtha depropanizer overhead, C_3-rich acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked hydrocarbons and treated to remove acidic impurities. It consists of hydrocarbons having carbon numbers in the range of C_2 through C_4, predominantly C_3.]	270-755-0	68477-73-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-063-00-1	Gases (petroleum), catalytic cracker; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic cracking process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-756-6	68477-74-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-064-00-7	Gases (petroleum), catalytic cracker, C_{1-5}-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_6, predominantly C_1 through C_5.]	270-757-1	68477-75-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-065-00-2	Gases (petroleum), catalytic polymd. naphtha stabilizer overhead, C_{2-4}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of catalytic polymerized naphtha. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_2 through C_6, predominantly C_2 through C_4.]	270-758-7	68477-76-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-066-00-8	Gases (petroleum), catalytic reformer, C_{1-4}-rich; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers in the range of C_1 through C_6, predominantly C_1 through C_4.]	270-760-8	68477-79-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-067-00-3	Gases (petroleum), C_{3-5} olefinic-paraffinic alkylation feed; Petroleum gas; [A complex combination of olefinic and paraffinic hydrocarbons having carbon numbers in the range of C_3 through C_5 which are used as alkylation feed. Ambient temperatures normally exceed the critical temperature of these combinations.]	270-765-5	68477-83-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-068-00-9	Gases (petroleum), C_4-rich; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from a catalytic fractionation process. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_4.]	270-767-6	68477-85-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-069-00-4	Gases (petroleum), deethanizer overheads; Petroleum gas; [A complex combination of hydrocarbons produced from distillation of the gas and gasoline fractions from the catalytic cracking process. It contains predominantly ethane and ethylene.]	270-768-1	68477-86-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-070-00-X	Gases (petroleum), deisobutanizer tower overheads; Petroleum gas; [A complex combination of hydrocarbons produced by the atmospheric distillation of a butane-butylene stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	270-769-7	68477-87-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-071-00-5	Gases (petroleum), depropanizer dry, propene-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from the gas and gasoline fractions of a catalytic cracking process. It consists predominantly of propylene with some ethane and propane.]	270-772-3	68477-90-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-072-00-0	Gases (petroleum), depropanizer overheads; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from the gas and gasoline fractions of a catalytic cracking process. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	270-773-9	68477-91-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-073-00-6	Gases (petroleum), gas recovery plant depropanizer overheads; Petroleum gas; [A complex combination of hydrocarbons obtained by fractionation of miscellaneous hydrocarbon streams. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4, predominantly propane.]	270-777-0	68477-94-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-074-00-1	Gases (petroleum), Girbatol unit feed; Petroleum gas; [A complex combination of hydrocarbons that is used as the feed into the Girbatol unit to remove hydrogen sulfide. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	270-778-6	68477-95-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-075-00-7	Gases (petroleum), isomerized naphtha fractionator, C_4-rich, hydrogen sulfide-free; Petroleum gas	270-782-8	68477-99-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-076-00-2	Tail gas (petroleum), catalytic cracked clarified oil and thermal cracked vacuum residue fractionation reflux drum; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked clarified oil and thermal cracked vacuum residue. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-802-5	68478-21-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-077-00-8	Tail gas (petroleum), catalytic cracked naphtha stabilization absorber; Petroleum gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-803-0	68478-22-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-078-00-3	Tail gas (petroleum), catalytic cracker, catalytic reformer and hydrodesulfurizer combined fractionater; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of products from catalytic cracking, catalytic reforming and hydrodesulfurizing processes treated to remove acidic impurities. It consists predominantly of hydrocarbons having cabon numbers predominantly in the range of C_1 through C_5.]	270-804-6	68478-24-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-079-00-9	Tail gas (petroleum), catalytic reformed naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of catalytic reformed naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	270-806-7	68478-26-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-080-00-4	Tail gas (petroleum), saturate gas plant mixed stream, C_4-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of straight-run naphtha, distillation tail gas and catalytic reformed naphtha stabilizer tail gas. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly butane and isobutane.]	270-813-5	68478-32-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-081-00-X	Tail gas (petroleum), saturate gas recovery plant, C_{1-2}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of distillate tail gas, straight-run naphtha, catalytic reformed naphtha stabilizer tail gas. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1through C_5, predominantly methane and ethane.]	270-814-0	68478-33-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-082-00-5	Tail gas (petroleum), vacuum residues thermal cracker; Petroleum gas; [A complex combination of hydrocarbons obtained from the thermal cracking of vacuum residues. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-815-6	68478-34-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-083-00-0	Hydrocarbons, C_{3-4}-rich, petroleum distillate; Petroleum gas; [A complex combination of hydrocarbons produced by distillation and condensation of crude oil. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_3 through C_4.]	270-990-9	68512-91-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-084-00-6	Gases (petroleum), full-range straight-run naphtha dehexanizer off; petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of the full-range straight-run naphtha. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-000-8	68513-15-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-085-00-1	Gases (petroleum), hydrocracking depropanizer off, hydrocarbon-rich; Petroleum gas; [A complex combination of hydrocarbon produced by the distillation of products from a hydrocracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4. It may also contain small amounts of hydrogen and hydrogen sulfide.]	271-001-3	68513-16-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-086-00-7	Gases (petroleum), light straight-run naphtha stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the stabilization of light straight-run naphtha. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-002-9	68513-17-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-087-00-2	Residues (petroleum), alkylation splitter, C_4-rich; Petroleum gas; [A complex residuum from the distillation of streams various refinery operations. It consists of hydrocarbons having carbon numbers in the range of C_4 through C_5, predominantly butane and boiling in the range of approximately - 11.7 °C to 27.8 °C (11 °F to 82 °F).]	271-010-2	68513-66-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-088-00-8	Hydrocarbons, C_{1-4}; Petroleum gas; [A complex combination of hydrocarbons provided by thermal cracking and absorber operations and by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately minus 164 °C to minus 0.5 °C (- 263 °F to 31 °F).]	271-032-2	68514-31-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-089-00-3	Hydrocarbons, C_{1-4}, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting hydrocarbon gases to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately - 164 °C to - 0.5 °C (-263 °F to 31 °F).]	271-038-5	68514-36-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-090-00-9	Hydrocarbons, C_{1-3}; Petroleum gas; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3 and boiling in the range of approximately minus 164 °C to minus 42 °C (- 263 °F to - 44 °F).]	271-259-7	68527-16-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-091-00-4	Hydrocarbons, C_{1-4}, debutanizer fraction; Petroleum gas	271-261-8	68527-19-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-092-00-X	Gases (petroleum), C_{1-5}, wet; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of crude oil and/or the cracking of tower gas oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	271-624-0	68602-83-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-093-00-5	Hydrocarbons, C_{2-4}; Petroleum gas	271-734-9	68606-25-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-094-00-0	Hydrocarbons, C_3; Petroleum gas	271-735-4	68606-26-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-095-00-6	Gases (petroleum), alkylation feed; Petroleum gas; [A complex combination of hydrocarbons produced by the catalytic cracking of gas oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	271-737-5	68606-27-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-096-00-1	Gases (petroleum), depropanizer bottoms fractionation off; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of depropanizer bottoms. It consists predominantly of butane, isobutane and butadiene.]	271-742-2	68606-34-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-097-00-7	Gases (petroleum), refinery blend; Petroleum gas; [A complex combination obtained from various processes. It consists of hydrogen, hydrogen sulfide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-183-7	68783-07-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-098-00-2	Gases (petroleum), catalytic cracking; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5.]	272-203-4	68783-64-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-099-00-8	Gases (petroleum), C_{2-4}, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4 and boiling in the range of approximately - 51 °C to - 34 °C (-60 °F to - 30 °F).]	272-205-5	68783-65-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-100-00-1	Gases (petroleum), crude oil fractionation off; Petroleum gas; [A complex combination of hydrocarbons produced by the fractionation of crude oil. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-871-7	68918-99-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-101-00-7	Gases (petroleum), dehexanizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of combined naphtha streams. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-872-2	68919-00-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-102-00-2	Gases (petroleum), light straight run gasoline fractionation stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of light straight-run gasoline. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-878-5	68919-05-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-103-00-8	Gases (petroleum), naphtha unifiner desulfurization stripper off; Petroleum gas; [A complex combination of hydrocarbons produced by a naphtha unifiner desulfurization process and stripped from the naphtha product. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-879-0	68919-06-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-104-00-3	Gases (petroleum), straight-run naphtha catalytic reforming off; Petroleum gas; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha and fractionation of the total effluent. It consists of methane, ethane, and propane.]	272-882-7	68919-09-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-105-00-9	Gases (petroleum), fluidized catalytic cracker splitter overheads; Petroleum gas; [A complex combination of hydrocarbons produced by the fractionation of the charge to the C_3 -C_4 splitter. It consists predominantly of C_3 hydrocarbons.]	272-893-7	68919-20-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-106-00-4	Gases (petroleum), straight-run stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of the liquid from the first tower used in the distillation of crude oil. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-883-2	68919-10-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-107-00-X	Gases (petroleum), catalytic cracked naphtha debutanizer; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked naphtha. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-169-3	68952-76-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-108-00-5	Tail gas (petroleum), catalytic cracked distillate and naphtha stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of catalytic cracked naphtha and distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-170-9	68952-77-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-109-00-0	Tail gas (petroleum), thermal-cracked distillate, gas oil and naphtha absorber; petroleum gas; [A complex combination of hydrocarbons obtained from the separation of thermal-cracked distillates, naphtha and gas oil. It consists pedrominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-175-6	68952-81-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-110-00-6	Tail gas (petroleum), thermal cracked hydrocarbon fractionation stabilizer, petroleum coking; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of thermal cracked hydrocarbons from petroleum coking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-176-1	68952-82-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-111-00-1	Gases (petroleum, light steam-cracked, butadiene conc.; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a thermal cracking process, It consists of hydrocarbons having a carbon number predominantly of C_4.]	273-265-5	68955-28-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-112-00-7	Gases (petroleum), straight-run naphtha catalytic reformer stabilizer overhead; Petroleum gas; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha and the fractionation of the total effluent. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	273-270-2	68955-34-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-113-00-2	Hydrocarbons, C_4; Petroleum gas	289-339-5	87741-01-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-114-00-8	Alkanes, C_{1-4}, C_3-rich; Petroleum gas	292-456-4	90622-55-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-115-00-3	Gases (petroleum), steam-cracker C_3-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a steam cracking process. It consists predominantly of propylene with some propane and boils in the range of approximately - 70 °C to 0 °C (- 94 °F to 32 °F).]	295-404-9	92045-22-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-116-00-9	Hydrocarbons, C_4, steam-cracker distillate; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products of a steam cracking process. It consists predominantly of hydrocarbons having a carbon number of C_4, predominantly 1-butene and 2-butene, containing also butane and isobutene and boiling in the range of approximately minus 12 °C to 5 °C (10.4 °F to 41 °F).]	295-405-4	92045-23-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-117-00-4	Petroleum gases, liquefied, sweetened, C_4 fraction; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting a liquified petroleum gas mix to a sweetening process to oxidize mercaptans or to remove acidic impurities. It consists predominantly of C_4 saturated and unsaturated hydrocarbons.]	295-463-0	92045-80-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			HKSU
649-118-00-X	Hydrocarbons, C_4, 1,3-butadiene- and isobutene-free; Petroleum gas	306-004-1	95465-89-7	Flam. Gas 1 Press. Gas Carc. 1B	H220 H350	GHS02 GHS04 GHS08 Dgr	H220 H350			H K U
649-119-00-5	Raffinates (petroleum), steam-cracked C_4 fraction cuprous ammonium acetate extn., C_{3-5} and C_{3-5} unsatd., butadiene-free; Petroleum gas	307-769-4	97722-19-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-120-00-0	Gases (petroleum), amine system feed; Refinery gas; [The feed gas to the amine system for removal of hydrogen sulfide. It consists of hydrogen. Carbon monoxide, carbon dioxide, hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5 may also be present.]	270-746-1	68477-65-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-121-00-6	Gases (petroleum), benzene unit hydrodesulfurizer off; Refinery gas; [Off gases produced by the benzene unit. It consists primarily of hydrogen. Carbon monoxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6, including benzene, may also be present.]	270-747-7	68477-66-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-122-00-1	Gases (petroleum), benzene unit recycle, hydrogen-rich; Refinery gas; [A complex combination of hydrocarbons obtained by recycling the gases of the benzene unit. It consists primarily of hydrogen with various small amounts of carbon monoxide and hydrocarbons having carbon numbers in the range of C_1 through C_6.]	270-748-2	68477-67-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-123-00-7	Gases (petroleum), blend oil, hydrogen-nitrogen-rich; Refinery gas; [A complex combination of hydrocarbons obtained by distillation of a blend oil. It consists primarily of hydrogen and nitrogen with various small amounts of carbon monoxide, carbon dioxide, and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-749-8	68477-68-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-124-00-2	Gases (petroleum), catalytic reformed naphtha stripper overheads; Refinery gas; [A complex combination of hydrocarbons obtained from stabilization of catalytic reformed naphtha. Its consists of hydrogen and saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	270-759-2	68477-77-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-125-00-8	Gases (petroleum), C_{6-8} catalytic reformer recycle; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from catalytic reforming of C_6-C_8 feed and recycled to conserve hydrogen. It consists primarily of hydrogen. It may also contain various small amounts of carbon monoxide, carbon dioxide, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-761-3	68477-80-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-126-00-3	Gases (petroleum), C_{6-8} catalytic reformer; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from catalytic reforming of C_6-C_8feed. It consists of hydrocarbons having carbon numbers in the range of C_1 through C_5 and hydrogen.]	270-762-9	68477-81-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-127-00-9	Gases (petroleum), C_{6-8} catalytic reformer recycle, hydrogen-rich; Refinery gas	270-763-4	68477-82-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-128-00-4	Gases (petroleum), C_2-return stream; Refinery gas; [A complex combination of hydrocarbons obtained by the extraction of hydrogen from a gas stream which consists primarily of hydrogen with small amounts of nitrogen, carbon monoxide, methane, ethane, and ethylene. It contains predominantly hydrocarbons such as methane, ethane, and ethylene with small amounts of hydrogen, nitrogen and carbon monoxide.]	270-766-0	68477-84-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-129-00-X	Gases (petroleum), dry sour, gas-concn.-unit-off; Refinery gas; [The complex combination of dry gases from a gas concentration unit. It consists of hydrogen, hydrogen sulfide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	270-774-4	68477-92-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-130-00-5	Gases (petroleum), gas concn. reabsorber distn.; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from combined gas streams in a gas concentration reabsorber. It consists predominantly of hydrogen, carbon monoxide, carbon dioxide, nitrogen, hydrogen sulfide and hydrocarbons having carbon numbers in the range of C_1 through C_3.]	270-776-5	68477-93-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-131-00-0	Gases (petroleum), hydrogen absorber off; Refinery gas; [A complex combination obtained by absorbing hydrogen from a hydrogen rich stream. It consists of hydrogen, carbon monoxide, nitrogen, and methane with small amounts of C_2 hydrocarbons.]	270-779-1	68477-96-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-132-00-6	Gases (petroleum), hydrogen-rich; Refinery gas; [A complex combination separated as a gas from hydrocarbon gases by chilling. It consists primarily of hydrogen with various small amounts of carbon monoxide, nitrogen, methane, and C_2 hydrocarbons.]	270-780-7	68477-97-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-133-00-1	Gases (petroleum), hydrotreater blend oil recycle, hydrogen-nitrogen-rich; Refinery gas; [A complex combination obtained from recycled hydrotreated blend oil. It consists primarily of hydrogen and nitrogen with various small amounts of carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-781-2	68477-98-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-134-00-7	Gases (petroleum), recycle, hydrogen-rich; Refinery gas; [A complex combination obtained from recycled reactor gases. It consists primarily of hydrogen with various small amounts of carbon monoxide, carbon dioxide, nitrogen, hydrogen sulfide, and saturated aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_5.]	270-783-3	68478-00-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-135-00-2	Gases (petroleum), reformer make-up, hydrogen-rich; Refinery gas; [A complex combination obtained from the reformers. It consists primarily of hydrogen with various small amounts of carbon monoxide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-784-9	68478-01-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-136-00-8	Gases (petroleum), reforming hydrotreater; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen, methane, and ethane with various small amounts of hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 thorugh C_5.]	270-785-4	68478-02-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-137-00-3	Gases (petroleum), reforming hydrotreater, hydrogen-methane-rich; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen and methane with various small amounts of carbon monoxide, carbon dioxide, nitrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_5.]	270-787-5	68478-03-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-138-00-9	Gases (petroleum), reforming hydrotreater make-up, hydrogen-rich; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen with various small amounts of carbon monoxide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-788-0	68478-04-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-139-00-4	Gases (petroleum), thermal cracking distn.; Refinery gas; [A complex combination produced by distillation of products from a thermal cracking process. It consists of hydrogen, hydrogen sulfide, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-789-6	68478-05-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-140-00-X	Tail gas (petroleum), catalytic cracker refractionation absorber; Refinery gas; [A complex combination of hydrocarbons obtained from refractionation of products from a catalytic cracking process. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	270-805-1	68478-25-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-141-00-5	Tail gas (petroleum), catalytic reformed naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from the catalytic reforming of straight run naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-807-2	68478-27-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-142-00-0	Tail gas (petroleum), catalytic reformed naphtha stabilizer; Refinery gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic reformed naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-808-8	68478-28-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-143-00-6	Tail gas (petroleum), cracked distillate hydrotreater separator; Refinery gas; [A complex combination of hydrocarbons obtained by treating cracked distillates with hydrogen in the presence of a catalyst. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-809-3	68478-29-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-144-00-1	Tail gas (petroleum), hydrodesulfurized straight-run naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from hydrodesulfurization of straight-run naphtha. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-810-9	68478-30-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-145-00-7	Gases (petroleum), catalytic reformed straight-run naphtha stabilizer overheads; Refinery gas; [A complex combination of hydrocarbons obtained from the catalytic reforming of straight-run naphtha followed by fractionation of the total effluent. It consists of hydrogen, methane, ethane and propane.]	270-999-8	68513-14-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-146-00-2	Gases (petroleum), reformer effluent high-pressure flash drum off; Refinery gas; [A complex combination produced by the high-pressure flashing of the effluent from the reforming reactor. It consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-003-4	68513-18-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-147-00-8	Gases (petroleum), reformer effluent low-pressure flash drum off; Refinery gas; [A complex combination produced by low-pressure flashing of the effluent from the reforming reactor. It consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-005-5	68513-19-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-148-00-3	Gases (petroleum), oil refinery gas distn. off; Refinery gas; [A complex combination separated by distillation of a gas stream containing hydrogen, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers in the range of C_1 through C_6 or obtained by cracking ethane and propane. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_2, hydrogen, nitrogen, and carbon monoxide.]	271-258-1	68527-15-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-149-00-9	Gases (petroleum), benzene unit hydrotreater depentanizer overheads; Refinery gas; [A complex combination produced by treating the feed from the benzene unit with hydrogen in the presence of a catalyst followed by depentanizing. It consists primarily of hydrogen, ethane and propane with various small amounts of nitrogen, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6. It may contain trace amounts of benzene.]	271-623-5	68602-82-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-150-00-4	Gases (petroleum), secondary absorber off, fluidized catalytic cracker overheads fractionator; Refinery gas; [A complex combination produced by the fractionation of the overhead products from the catalytic cracking process in the fluidized catalytic cracker. It consists of hydrogen, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	271-625-6	68602-84-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-151-00-X	Petroleum products, refinery gases; Refinery gas; [A complex combination which consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-750-6	68607-11-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-152-00-5	Gases (petroleum), hydrocracking low-pressure separator; Refinery gas; [A complex combination obtained by the liquid-vapor separation of the hydrocracking process reactor effluent. It consists predominantly of hydrogen and saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	272-182-1	68783-06-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-153-00-0	Gases (petroleum), refinery; Refinery gas; [A complex combination obtained from various petroleum refining operations. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	272-338-9	68814-67-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-154-00-6	Gases (petroleum), platformer products separator off; Refinery gas; [A complex combination obtained from the chemical reforming of naphthenes to aromatics. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	272-343-6	68814-90-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-155-00-1	Gases (petroleum), hydrotreated sour kerosine depentanizer stabilizer off; Refinery gas; [The complex combination obtained from the depentanizer stabilization of hydrotreated kerosine. It consists primarily of hydrogen, methane, ethane, and propane with various small amounts of nitrogen, hydrogen sulfide, carbon monoxide and hydrocarbons having carbon numbers predominantly in the range of C_4 through C_5.]	272-775-5	68911-58-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-156-00-7	Gases (petroleum), hydrotreated sour kerosine flash drum; Refinery gas; [A complex combination obtained from the flash drum of the unit treating sour kerosine with hydrogen in the presence of a catalyst. It consists primarily of hydrogen and methane with various small amounts of nitrogen, carbon monoxide, and hydrocarbons having carbon numbers predominantly in the range of C_2 through C_5.]	272-776-0	68911-59-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-157-00-2	Gases (petroleum), distillate unifiner desulfurization stripper off; Refinery gas; [A complex combination stripped from the liquid product of the unifiner desulfurization process. It consists of hydrogen sulfide, methane, ethane, and propane.]	272-873-8	68919-01-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-158-00-8	Gases (petroleum), fluidized catalytic cracker fractionation off; Refinery gas; [A complex combination produced by the fractionation of the overhead product of the fluidized catalytic cracking process. It consists of hydrogen, hydrogen sulfide, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-874-3	68919-02-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-159-00-3	Gases (petroleum), fluidized catalytic cracker scrubbing secondary absorber off; Refinery gas; [A complex combination produced by scrubbing the overhead gas from the fluidized catalytic cracker. It consists of hydrogen, nitrogen, methane, ethane and propane.]	272-875-9	68919-03-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-160-00-9	Gases (petroleum), heavy distillate hydrotreater desulfurization stripper off; Refinery gas; [A complex combination stripped from the liquid product of the heavy distillate hydrotreater desulfurization process. It consists of hydrogen, hydrogen sulfide, and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5].	272-876-4	68919-04-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-161-00-4	Gases (petroleum), platformer stabilizer off, light ends fractionation; Refinery gas; [A complex combination obtained by the fractionation of the light ends of the platinum reactors of the plattformer unit. It consists of hydrogen, methane, ethane and propane.]	272-880-6	68919-07-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-162-00-X	Gases (petroleum), preflash tower off, crude distn.; Refinery gas; [A complex combination produced from the first tower used in the distillation of crude oil. It consists of nitrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-881-1	68919-08-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-163-00-5	Gases (petroleum), tar stripper off; Refinery gas; [A complex combination obtained by the fractionation of reduced crude oil. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-884-8	68919-11-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-164-00-0	Gases (petroleum), unifiner stripper off; Refinery gas; [A combination of hydrogen and methane obtained by fractionation of the products from the unifiner unit.]	272-885-3	68919-12-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-165-00-6	Tail gas (petroleum), catalytic hydrodesulfurized naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from the hydrodesulfurization of naphtha. It consists of hydrogen, methane, ethane, and propane.]	273-173-5	68952-79-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-166-00-1	Tail gas (petroleum), straight-run naphtha hydrodesulfurizer; Refinery gas; [A complex combination obtained from the hydrodesulfurization of straight-run naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	273-174-0	68952-80-7	Press. Gas Flam. Gas 1 Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-167-00-7	Gases (petroleum), sponge absorber off, fluidized catalytic cracker and gas oil desulfurizer overhead fractionation; Refinery gas; [A complex combination obtained by the fractionation of products from the fluidized catalytic cracker and gas oil desulfurizer. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-269-7	68955-33-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-168-00-2	Gases (petroleum), crude distn. and catalytic cracking; Refinery gas; [A complex combination produced by crude distillation and catalytic cracking processes. It consists of hydrogen, hydrogen sulfide, nitrogen, carbon monoxide and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-563-5	68989-88-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-169-00-8	Gases (petroleum), gas oil diethanolamine scrubber off; Refinery gas; [A complex combination produced by desulfurization of gas oils with diethanolamine. It consists predominantly of hydrogen sulfide, hydrogen and aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_5.]	295-397-2	92045-15-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-170-00-3	Gases (petroleum), gas oil hydrodesulfurization effluent; Refinery gas; [A complex combination obtained by separation of the liquid phase from the effluent from the hydrogenation reaction. It consists predominantly of hydrogen, hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	295-398-8	92045-16-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-171-00-9	Gases (petroleum), gas oil hydrodesulfurization purge; Refinery gas; [A complex combination of gases obtained from the reformer and from the purges from the hydrogenation reactor. It consists predominantly of hydrogen and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	295-399-3	92045-17-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-172-00-4	Gases (petroleum), hydrogenator effluent flash drum off; Refinery gas; [A complex combination of gases obtained from flash of the effluents after the hydrogenation reaction. It consists predominantly of hydrogen and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	295-400-7	92045-18-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-173-00-X	Gases (petroleum), naphtha steam cracking high-pressure residual; Refinery gas; [A complex combination obtained as a mixture of the non-condensable portions from the product of a naphtha steam cracking process as well as residual gases obtained during the preparation of subsequent products. It consists predominantly of hydrogen and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5 with which natural gas may also be mixed.]	295-401-2	92045-19-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-174-00-5	Gases (petroleum), residue visbaking off; Refinery gas; [A complex combination obtained from viscosity reduction of residues in a furnace. It consists predominantly of hydrogen sulfide and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	295-402-8	92045-20-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-175-00-0	Foots oil (petroleum), acid-treated; Foots oil; [A complex combination of hydrocarbons obtained by treatment of Foot's oil with sulfuric acid. It consists predominantly of branched-chain hydrocarbons with carbon numbers predominantly in the range of C_{20} through C_{50}.]	300-225-7	93924-31-3	Flam. Gas 1 Press. Gas Carc. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-176-00-6	Foots oil (petroleum), clay-treated; Foots oil; [A complex combination of hydrocarbons obtained by treatment of Foot's oil with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists predominantly of branched chain hydrocarbons with carbon numbers predominantly in the range of C_{20} through C_{50}.]	300-226-2	93924-32-4	Flam. Gas 1 Press. Gas Carc. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-177-00-1	Gases (petroleum), C_{3-4}; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from the cracking of crude oil. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_4, predominantly of propane and propylene, and boiling in the range of approximately - 51 °C to - 1 °C (-60°F to 30 °F.)]	268-629-5	68131-75-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-178-00-7	Tail gas (petroleum), catalytic cracked distillate and catalytic cracked naphtha fractionation absorber; Petroleum gas; [The complex combination of hydrocarbons from the distillation of the products from catalytic cracked distillates and catalytic cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4.]	269-617-2	68307-98-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-179-00-2	Tail gas (petroleum), catalytic polymn. naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons from the fractionation stabilization products from polymerization of naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4.]	269-618-8	68307-99-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-180-00-8	Tail gas (petroleum), catalytic reformed naphtha fractionation stabilizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation stabilization of catalytic reformed naphtha and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-619-3	68308-00-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-181-00-3	Tail gas (petroleum), cracked distillate hydrotreater stripper; Petroleum gas; [A complex combination of hydrocarbons obtained by treating thermal cracked distillates with hydrogen in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-620-9	68308-01-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-182-00-9	Tail gas (petroleum), straight-run distillate hydrodesulfurizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from catalytic hydrodesulfurization of straight run distillates and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-630-3	68308-10-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-183-00-4	Tail gas (petroleum), gas oil catalytic cracking absorber; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of products from the catalytic cracking of gas oil. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-623-5	68308-03-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-184-00-X	Tail gas (petroleum), gas recovery plant; Petroleum gas; [A complex combination of hydrocarbons from the distillation of products from miscellaneous hydrocarbon streams. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-624-0	68308-04-3	Press. Gas Flam. Gas 1 Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-185-00-5	Tail gas (petroleum), gas recovery plant deethanizer; Petroleum gas; [A complex combination of hydrocarbons from the distillation of products from miscellaneous hydrocarbon streams. It consists of hydrocarbon having carbon numbers predominantly in the range of C_1 through C_4.]	269-625-6	68308-05-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-186-00-0	Tail gas (petroleum), hydrodesulfurized distillate and hydrodesulfurized naphtha fractionator, acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of hydrodesulfurized naphtha and distillate hydrocarbon streams and treated to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-626-1	68308-06-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-187-00-6	Tail gas (petroleum), hydrodesulfurized vacuum gas oil stripper, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from stripping stabilization of catalytic hydrodesulfurized vacuum gas oil and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-627-7	68308-07-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-188-00-1	Tail gas (petroleum), light straight-run naphtha stabilizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation stabilization of light straight run naphtha and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-629-8	68308-09-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-189-00-7	Tail gas (petroleum), propane-propylene alkylation feed prep deethanizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of the reaction products of propane with propylene. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-631-9	68308-11-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-190-00-2	Tail gas (petroleum), vacuum gas oil hydrodesulfurizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from catalytic hydrodesulfurization of vacuum gas oil and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-632-4	68308-12-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-191-00-8	Gases (petroleum), catalytic cracked overheads; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from the catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5 and boiling in the range of approximately - 48 °C to 32 °C (- 54 °F to 90 °F).]	270-071-2	68409-99-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-193-00-9	Alkanes, C_{1-2}; Petroleum gas	270-651-5	68475-57-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-194-00-4	Alkanes, C_{2-3}; Petroleum gas	270-652-0	68475-58-1	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-195-00-X	Alkanes, C_{3-4}; Petroleum gas	270-653-6	68475-59-2	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-196-00-5	Alkanes, C_{4-5}; Petroleum gas	270-654-1	68475-60-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-197-00-0	Fuel gases; Petroleum gas; [A combination of light gases. It consists predominantly of hydrogen and/or low molecular weight hydrocarbons.]	270-667-2	68476-26-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-198-00-6	Fuel gases, crude oil of distillates; Petroleum gas; [A complex combination of light gases produced by distillation of crude oil and by catalytic reforming of naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately - 217 °C to - 12 °C (- 423 °F to 10 °F).]	270-670-9	68476-29-9	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-199-00-1	Hydrocarbons, C_{3-4}; Petroleum gas	270-681-9	68476-40-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-200-00-5	Hydrocarbons, C_{4-5}; Petroleum gas	270-682-4	68476-42-6	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-201-00-0	Hydrocarbons, C_{2-4}, C_3-rich; Petroleum gas	270-689-2	68476-49-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-202-00-6	Petroleum gases, liquefied; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_7 and boiling in the range of approximately - 40 °C to 80 °C (- 40 °F to 176 °F).]	270-704-2	68476-85-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			HKSU

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649-203-00-1	Petroleum gases, liquefied, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting liquefied petroleum gas mix to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_7 and boiling in the range of approximately - 40 °C to 80 °C (- 40 °F to 176 °F).]	270-705-8	68476-86-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			HKSU
649-204-00-7	Gases (petroleum), C_{3-4}, isobutane-rich; Petroleum gas; [A complex combination of hydrocarbons from the distillation of saturated and unsaturated hydrocarbons usually ranging in carbon numbers from C_3 through C_6, predominantly butane and isobutane. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_4, predominantly isobutane.]	270-724-1	68477-33-8	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-205-00-2	Distillates (petroleum), C_{3-6}, piperylene-rich; Petroleum gas; [A complex combination of hydrocarbons from the distillation of saturated and unsaturated aliphatic hydrocarbons usually ranging in the carbon numbers C_3 through C_6. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly piperylenes.]	270-726-2	68477-35-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-206-00-8	Gases (petroleum), butane splitter overheads; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of the butane stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	270-750-3	68477-69-0	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-207-00-3	Gases (petroleum), C_{2-}; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic fractionation process. It contains predominantly ethane, ethylene, propane, and propylene.]	270-751-9	68477-70-3	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-208-00-9	Gases (petroleum), catalytic-cracked gas oil depropanizer bottoms, C_4-rich acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked gas oil hydrocarbon stream and treated to remove hydrogen sulfide and other acidic components. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_4.]	270-752-4	68477-71-4	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-209-00-4	Gases (petroleum), catalytic-cracked naphtha debutanizer bottoms, C_{3-5}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic cracked naphtha. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5.]	270-754-5	68477-72-5	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U
649-210-00-X	Tail gas (petroleum), isomerized naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization products from isomerized naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-628-2	68308-08-7	Flam. Gas 1 Press. Gas Carc. 1A Muta. 1B	H220 H350 H340	GHS02 GHS04 GHS08 Dgr	H220 H350 H340			H K U

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649-211-00-5	Foots oil (petroleum), carbon-treated; Foots oil; [A complex combination of hydrocarbons obtained by the treatment of Foots oil with activated carbon for the removal of trace constituents and impurities. It consists predominantly of saturated straight chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-126-0	97862-76-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-212-00-0	Distillates (petroleum), sweetened middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-088-7	64741-86-2	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-213-00-6	Gas oils (petroleum), solvent-refined; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-092-9	64741-90-8	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-214-00-1	Distillates (petroleum), solvent-refined middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-093-4	64741-91-9	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-215-00-7	Gas oils (petroleum), acid-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-112-6	64742-12-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-216-00-2	Distillates (petroleum), acid-treated middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{20} and boiling in the range of approximately 205 °C to 345 °C (401 °F to 653 °F).]	265-113-1	64742-13-8	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-217-00-8	Distillates (petroleum), acid-treated light; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{9} through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-114-7	64742-14-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-218-00-3	Gas oils (petroleum), chemically neutralized; Gasoil — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-129-9	64742-29-6	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-219-00-9	Distillates (petroleum), chemically neutralized middle; Gasoil — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{20} and boiling in the range of approximately 205 °C to 345 °C (401 °F to 653 °F).]	265-130-4	64742-30-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-220-00-4	Distillates (petroleum), clay-treated middle; Gasoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-139-3	64742-38-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-221-00-X	Distillates (petroleum), hydrotreated middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-148-2	64742-46-7	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-222-00-5	Gas oils (petroleum), hydrodesulfurized; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-182-8	64742-79-6	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-223-00-0	Distillates (petroleum), hydrodesulfurized middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-183-3	64742-80-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-224-00-6	Fuels, diesel; Gasoil — unspecified; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 163 °C to 357 °C (325 °F to 675 °F).]	269-822-7	68334-30-5	Carc. 2	H351	GHS08 Wng	H351			H N
649-225-00-1	Fuel oil, No 2; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 32,6 SUS at 37,7 °C (100 °F) to a maximum of 37,9 SUS at 37,7 °C (100 °F).]	270-671-4	68476-30-2	Carc. 2	H351	GHS08 Wng	H351			H

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649-226-00-7	Fuel oil, No 4; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 45 SUS at 37,7 °C (100 °F) to a maximum of 125 SUS at 37,7 °C (100 °F).]	270-673-5	68476-31-3	Carc. 2	H351	GHS08 Wng	H351			H
649-227-00-2	Fuels, diesel, No 2; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 32,6 SUS at 37,7 °C (100 °F).]	270-676-1	68476-34-6	Carc. 2	H351	GHS08 Wng	H351			H
649-228-00-8	Distillates (petroleum), catalytic reformer fractionator residue, high-boiling; Gasoil — unspecified; [A complex combination of hydrocarbons from the distillation of catalytic reformer fracftionator residue. It boils in the range of approximately 343 °C to 399 °C (650 °F to 750 °F).]	270-719-4	68477-29-2	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-229-00-3	Distillates (petroleum), catalytic reformer fractionator residue, intermediate-boiling; Gasoil — unspecified; [A complex combination of hydrocarbons from the distillation of catalytic reformer fractionator residue. It boils in the range of approximately 288 °C to 371 °C (550 °F to 700 °F).]	270-721-5	68477-30-5	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-230-00-9	Distillates (petroleum), catalytic reformer fractionator residue, low-boiling; Gasoil — unspecified; [The complex combination of hydrocarbons from the distillation of catalytic reformer fractionator residue. It boils approximately below 288 °C (550 °F).]	270-722-0	68477-31-6	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-231-00-4	Distillates (petroleum), highly refined middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the subjection of a petroleum fraction to several of the following steps: filtration, centrifugation, atmospheric distillation, vacuum distillation, acidification, neutralization and clay treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{20}.]	292-615-8	90640-93-0	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-232-00-X	Distillates (petroleum) catalytic reformer, heavy arom. conc.; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from the distillation of a catalytically reformed petroleum cut. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{16} and boiling in the range of approximately 200 °C to 300 °C (392 °F to 572 °F).]	295-294-2	91995-34-5	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-233-00-5	Gas oils, paraffinic; Gasoil — unspecified; [A distillate obtained from the redistillation of a complex combination of hydrocarbons obtained by the distillation of the effluents from a severe catalytic hydrotreatment of paraffins. It boils in the range of approximately 190 °C to 330 °C (374 °F to 594 °F).]	300-227-8	93924-33-5	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-234-00-0	Naphtha (petroleum), solvent-refined hydrodesulfurized heavy; Gasoil — unspecified	307-035-3	97488-96-5	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-235-00-6	Hydrocarbons, C_{16-20}, hydrotreated middle distillate, distn. lights; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of a middle distillate with hydrogen. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{20} and boiling in the range of approximately 290 °C to 350 °C (554 °F to 662 °F). It produces a finished oil having a viscosity of 2cSt at 100 °C (212 °F).]	307-659-6	97675-85-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-236-00-1	Hydrocarbons, C_{12-20}, hydrotreated paraffinic, distn. lights; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of heavy paraffins with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{20} and boiling in the range of approximately 230 °C to 350 °C (446 °F to 662 °F). It produces a finished oil having a viscosity of 2cSt at 100 °C (212 °F).]	307-660-1	97675-86-0	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-237-00-7	Hydrocarbons, C_{11-17}, solvent-extd. light naphthenic; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a visciosity of 2.2 cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{17} and boiling in the range of approximately 200 °C to 300 °C (392 °F to 572 °F).]	307-757-9	97722-08-2	Carc. 1B	H350	GHS08 Dgr	H350			H N

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649-238-00-2	Gas oils, hydrotreated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from the redistillation of the effluents from the treatment of paraffins with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{27} and boiling in the range of approximately 330 °C to 340 °C (626 °F to 644 °F).]	308-128-1	97862-78-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-239-00-8	Distillates (petroleum), carbon-treated light paraffinic; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of a petroleum oil fraction with activated charcoal for the removal of traces of polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{28}.]	309-667-5	100683-97-4	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-240-00-3	Distillates (petroleum), intermediate paraffinic, carbon-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of petroleum with activated charcoal for the removal of trace polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	309-668-0	100683-98-5	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-241-00-9	Distillates (petroleum), intermediate paraffinic, clay-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of petroleum with bleaching earth for the removal of trace polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	309-669-6	100683-99-6	Carc. 1B	H350	GHS08 Dgr	H350			H N

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-242-00-4	Alkanes, C_{12-26}-branched and linear	292-454-3	90622-53-0	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-243-00-X	Lubricating greases; Grease; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{50}. May contain organic salts of alkali metals, alkaline earth metals, and/or aluminium compounds.]	278-011-7	74869-21-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-244-00-5	Slack wax (petroleum); Slack wax; [A complex combination of hydrocarbons obtained from a petroleum fraction by solvent crystallization (solvent dewaxing) or as a distillation fraction from a very waxy crude. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	265-165-5	64742-61-6	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-245-00-0	Slack wax (petroleum), acid-treated; Slack wax; [A complex combination of hydrocarbons obtained as a raffinate by treatment of a petroleum slack wax fraction with sulfuric acid treating process. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	292-659-8	90669-77-5	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-246-00-6	Slack wax (petroleum), clay-treated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of a petroleum slack wax fraction with natural or modified clay in either a contacting or percolation process. It consists predominantly of saturated straight and branched hydrocarbons having carbon numbers predominantly greater than C_{20}.]	292-660-3	90669-78-6	Carc. 1B	H350	GHS08 Dgr	H350			H N

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-247-00-1	Slack wax (petroleum), hydrotreated; Slack wax; [A complex combination of hydrocarbons obtained by treating slack wax with hydrogen in the presence of a catalyst. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	295-523-6	92062-09-4	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-248-00-7	Slack wax (petroleum), low-melting; Slack wax; [A complex combination of hydrocarbons obtained from a petroleum fraction by solvent deparaffination. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-524-1	92062-10-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-249-00-2	Slack wax (petroleum), low-melting, hydrotreated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of low-melting petroleum slack wax with hydrogen in the presence of a catalyst. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-525-7	92062-11-8	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-250-00-8	Slack wax (petroleum), low-melting, carbon-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting slack wax with activated carbon for the removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-155-9	97863-04-2	Carc. 1B	H350	GHS08 Dgr	H350			H N

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-251-00-3	Slack wax (petroleum), low-melting, clay-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting petroleum slack wax with bentonite for removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-156-4	97863-05-3	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-252-00-9	Slack wax (petroleum), low-melting, silicic acid-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting petroleum slack wax with silicic acid for the removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-158-5	97863-06-4	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-253-00-4	Slack wax (petroleum), carbon-treated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of petroleum slack wax with activated charcoal for the removal of trace polar constituents and impurities.]	309-723-9	100684-49-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-254-00-X	Petrolatum; Petrolatum; [A complex combination of hydrocarbons obtained as a semi-solid from dewaxing paraffinic residual oil. It consists predominantly of saturated crystalline and liquid hydrocarbons having carbon numbers predominantly greater than C_{25}.]	232-373-2	8009-03-8	Carc. 1B	H350	GHS08 Dgr	H350			H N

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-255-00-5	Petrolatum (petroleum), oxidized; Petrolatum; [A complex combination of organic compounds, predominantly high molecular weight carboxylic acids, obtained by the air oxidation of petrolatum.]	265-206-7	64743-01-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-256-00-0	Petrolatum (petroleum), alumina-treated; Petrolatum; [A complex combination of hydrocarbons obtained when petrolatum is treated with Al_2O_3 to remove polar components and impurities. It consists predominantly of saturated, crystalline, and liquid hydrocarbons having carbon numbers predominantly greater than C_{25}.]	285-098-5	85029-74-9	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-257-00-6	Petrolatum (petroleum), hydrotreated; Petrolatum; [A complex combination of hydrocarbons obtained as a semi-solid from dewaxed paraffinic residual oil treated with hydrogen in the presence of a catalyst. It consists predominantly of saturated microcrystalline and liquid hydrocarbons having carbon numbers predominantly greater than C_{20}.]	295-459-9	92045-77-7	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-258-00-1	Petrolatum (petroleum), carbon-treated; Petrolatum; [A complex combination of hydrocarbons obtained by the treatment of petroleum petrolatum with activated carbon for the removal of trace polar constituents and impurities. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly greater than C_{20}.]	308-149-6	97862-97-0	Carc. 1B	H350	GHS08 Dgr	H350			H N

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-259-00-7	Petrolatum (petroleum), silicic acid-treated; Petrolatum; [A complex combination of hydrocarbons obtained by the treatment of petroleum petrolatum with silicic acid for the removal of trace polar constituents and impurities. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly greater than C_{20}.]	308-150-1	97862-98-1	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-260-00-2	Petrolatum (petroleum), clay-treated; Petrolatum; [A complex combination of hydrocarbons obtained by treatment of petrolatum with bleaching earth for the removal of traces of polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of greater than C_{25}.]	309-706-6	100684-33-1	Carc. 1B	H350	GHS08 Dgr	H350			H N
649-261-00-8	Gasoline, natural; Low boiling point naphtha; [A complex combination of hydrocarbons separated from natural gas by processes such as refrigeration or absorption. It consists predominantly of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_8 and boiling in the range of approximately minus 20 °C to 120 °C (- 4 °F to 248 °F).]	232-349-1	8006-61-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-262-00-3	Naphtha; Low boiling point naphtha; [Refined, partly refined, or unrefined petroleum products by the distillation of natural gas. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_6 and boiling in the range of approximately 100 °C to 200 °C (212 °F to 392 °F).]]	232-443-2	8030-30-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-263-00-9	Ligroine; Low boiling point naphtha; [A complex combination of hydrocarbons obtained by the fractional distillation of petroleum. This fraction boils in a range of approximately 20 °C to 135 °C (58 °F to 275 °F).]	232-453-7	8032-32-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-264-00-4	Naphtha (petroleum), heavy straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-041-0	64741-41-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-265-00-X	Naphtha (petroleum), full-range straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 220 °C (- 4 °F to 428 °F).]	265-042-6	64741-42-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-266-00-5	Naphtha (petroleum), light straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10} and boiling in the range of approximately minus 20 °C to 180 °C (- 4 °F to 356 °F).]	265-046-8	64741-46-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-267-00-0	Solvent naphtha (petroleum), light aliph.; Low boiling point naphtha; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{10} and boiling in the range of approximately 35 °C to 160 °C (95 °F to 320 °F).]	265-192-2	64742-89-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-268-00-6	Distillates (petroleum), straight-run light; Low boiling point naphtha; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_7 and boiling in the range of approximately - 88 °C to 99 °C (- 127 °F to 210 °F).]	270-077-5	68410-05-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-269-00-1	Gasoline, vapor-recovery; Low boiling point naphtha; [A complex combination of hydrocarbons separated from the gases from vapor recovery systems by cooling. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately - 20 °C to 196 °C (-4 °F to 384 °F).]	271-025-4	68514-15-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-270-00-7	Gasoline, straight-run, topping-plant; Low boiling point naphtha; [A complex combination of hydrocarbons produced from the topping plant by the distillation of crude oil. It boils in the range of approximately 36,1 °C to 193,3 °C (97 °F to 380 °F).]	271-727-0	68606-11-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-271-00-2	Naphtha (petroleum), unsweetened; Low boiling point naphtha; [A complex combination of hydrocarbons produced from the distillation of naphtha streams from various refinery processes. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 0 °C to 230 °C (25 °F to 446 °F).]	272-186-3	68783-12-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-272-00-8	Distillates (petroleum), light straight-run gasoline fractionation stabilizer overheads; Low boiling point naphtha; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_3 through $C_{6.}$]	272-931-2	68921-08-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-273-00-3	Naphtha (petroleum), heavy straight run, arom.-contg.; Low boiling point naphtha; [A complex combination of hydrocarbons obtained from a distillation process of crude petroleum. It consists predominantly of hydrocarbons having carbon numbers in the range of C_8 through C_{12} and boiling in the range of approximately 130 °C to 210 °C (266 °F to 410 °F).]	309-945-6	101631-20-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-274-00-9	Naphtha (petroleum), full-range alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consist of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 220 °C (194 °F to 428 °F).]	265-066-7	64741-64-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-275-00-4	Naphtha (petroleum), heavy alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 to C_5. It consists of predominantly branched chain saturated hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{12} and boiling in the range of approximately 150 °C to 220 °C (302 °F to 428 °F).]	265-067-2	64741-65-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-276-00-X	Naphtha (petroleum), light alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_7 through C_{10} and boiling in the range of aproximately 90 °C to 160 °C (194 °F to 320 °F).]	265-068-8	64741-66-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-277-00-5	Naphtha (petroleum), isomerization; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained from catalytic isomerization of straight chain paraffinic C_4 through C_6 hydrocarbons. It consists predominantly of saturated hydrocarbons such as isobutane, isopentane, 2,2-dimethylbutane, 2-methylpentane, and 3-methylpentane.]	265-073-5	64741-70-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-278-00-0	Naphtha (petroleum), solvent-refined light; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F).]	265-086-6	64741-84-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-279-00-6	Naphtha (petroleum), solvent-refined heavy; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-095-5	64741-92-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-280-00-1	Raffinates (petroleum), catalytic reformer ethylene glycol-water countercurrent exts.; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from the UDEX extraction process on the catalytic reformer stream. It consists of saturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_9.]	270-088-5	68410-71-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-281-00-7	Raffinates (petroleum), reformer, Lurgi unit-sepd.; Low boiling point modified naphtha; [The complex combination of hydrocarbons obtained as a raffinate from a Lurgi separation unit. It consists predominantly of non-aromatic hydrocarbons with various small amounts of aromatic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8.]	270-349-3	68425-35-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-282-00-2	Naphtha (petroleum), full-range alkylate, butane-contg.; Low boiling point modified naphta; [A complex combination of hydrocarbons produced by the distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} with some butanes and boiling in the range of approximately 35 °C to 200 °C (95 °F to 428 °F).]	271-267-0	68527-27-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-283-00-8	Distillates (petroleum), naphtha steam cracking-derived, solvent-refined light hydrotreated; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinates from a solvent extraction process of hydrotreated light distillate from steam-cracked naphtha.]	295-315-5	91995-53-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-284-00-3	Naphtha (petroleum), C_{4-12} butane-alkylate, isooctane-rich; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by alkylation of butanes. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12}, rich in isooctane, and boiling in the range of approximately 35 °C to 210 °C (95 °F to 410 °F).]	295-430-0	92045-49-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-285-00-9	Hydrocarbons, hydrotreated light naphtha distillates, solvent-refined; Low boiling point modified naphtha; [A combination of hydrocarbons obtained from the distillation of hydrotreated naphtha followed by a solvent extraction and distillation process. It consists predominantly of saturated hydrocarbons boiling in the range of approximately 94 °C to 99 °C (201 °F to 210 °F).]	295-436-3	92045-55-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-286-00-4	Naphtha (petroleum), isomerization, C_6-fraction; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by distillation of a gasoline which has been catalytically isomerized. It consists predominantly of hexane isomers boiling in the range of approximately 60 °C to 66 °C (140 °F to 151 °F).]	295-440-5	92045-58-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-287-00-X	Hydrocarbons, C_{6-7}, naphtha-cracking, solvent-refined; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by the sorption of benzene from a catalytically fully hydrogenated benzene-rich hydrocarbon cut that was distillatively obtained from prehydrogenated cracked naphtha. It consists predominantly of paraffinic and naphthenic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_7 and boiling in the range of approximately 70 °C to 100 °C (158 °F to 212 °F).]	295-446-8	92045-64-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-288-00-5	Hydrocarbons, C_6-rich, hydrotreated light naphtha distillates, solvent-refined; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by distillation of hydrotreated naphtha followed by solvent extraction. It consists predominantly of saturated hydrocarbons and boiling in the range of approximately 65 °C to 70 °C (149 °F to 158 °F).]	309-871-4	101316-67-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-289-00-0	Naphtha (petroleum), heavy catalytic cracked; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by a distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (148 °F to 446 °F). It contains a relatively large proportion of unsaturated hydrocarbons.]	265-055-7	64741-54-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-290-00-6	Naphtha (petroleum), light catalytic cracked; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F). It contains a relatively large proportion of unsaturated hydrocarbons.]	265-056-2	64741-55-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-291-00-1	Hydrocarbons, C_{3-11}, catalytic cracker distillates; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillations of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_{11} and boiling in a range approximately up to 204 °C (400 °F).]	270-686-6	68476-46-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-292-00-7	Naphtha (petroleum), catalytic cracked light distd.; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-185-8	68783-09-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-293-00-2	Distillates (petroleum), naphtha steam cracking-derived, hydrotreated light arom.; Low boiling point cat-cracked naphtha.; [A complex combination of hydrocarbons obtained by treating a light distillate from steam-cracked naphtha. It consists predominantly of aromatic hydrocarbons.]	295-311-3	91995-50-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-294-00-8	Naphtha (petroleum), heavy catalytic cracked, sweetened; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting a catalytic cracked petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 60 °C to 200 °C (140 °F to 392 °F).]	295-431-6	92045-50-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-295-00-3	Naphtha (petroleum), light catalytic cracked sweetened; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting naphtha from a catalytic cracking process to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons boiling in a range of approximately 35 °C to 210 °C (95 °F to 410 °F).]	295-441-0	92045-59-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-296-00-9	Hydrocarbons, C_{8-12}, catalytic-cracking, chem. neutralized; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of a cut from the catalytic cracking process, having undergone an alkaline washing. It consists predominantly of hydrocarbons having carbon numbers in the range of C_8 through C_{12} and boiling in the range of approximately 130 °C to 210 °C (266 °F to 410 °F).]	295-794-0	92128-94-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-297-00-4	Hydrocarbons, C_{8-12}, catalytic cracker distillates; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by distillation of products from a catalytic cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12} and boiling in the range of approximately 140 °C to 210 °C (284 °F to 410 °F).]	309-974-4	101794-97-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-298-00-X	Hydrocarbons, C_{8-12}, catalytic cracking, chem. neutralized, sweetened; Low boiling point cat-cracked naphtha	309-987-5	101896-28-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-299-00-5	Naphtha (petroleum), light catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced from the distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F). It contains a relatively large proportion of aromatic and branched chain hydrocarbons. This stream may contain 10 vol. % or more benzene.]]	265-065-1	64741-63-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-300-00-9	Naphtha (petroleum), heavy catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced from the distillation of products from a catalytic reforming process. It consists of predominantly aromatic hydrocarbons having numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-070-9	64741-68-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-301-00-4	Distillates (petroleum), catalytic reformed depentanizer; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons from the distillation of products from a catalytic reforming process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 49 °C to 63 °C - 57 °F to 145 °F).]	270-660-4	68475-79-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-302-00-X	Hydrocarbons, C_{2-6}, C_{6-8} catalytic reformer; Low boiling point cat-reformed naphtha	270-687-1	68476-47-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-303-00-5	Residues (petroleum), C_{6-8} catalytic reformer; Low boiling point cat-reformed naphtha'; [A complex residuum from the catalytic reforming of C_{6-8} feed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	270-794-3	68478-15-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-304-00-0	Naphtha (petroleum), light catalytic reformed, arom.-free; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained from distillation of products from a catalytic reforming process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_8 and boiling in the range of approximately 35 °C to 120 °C (95 °F to 248 °F). It contains a relatively large proportion of branched chain hydrocarbons with the aromatic components removed.]	270-993-5	68513-03-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-305-00-6	Distillates (petroleum), catalytic reformed straight-run naphtha overheads; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha followed by the fractionation of the total effluent. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-008-1	68513-63-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-306-00-1	Petroleum products, hydrofiner-powerformer reformates; Low boiling point cat-reformed naphtha; [The complex combination of hydrocarbons obtained in a hydrofiner-powerformer process and boiling in a range of approximately 27 °C to 210 °C (80 °F to 410 °F).]	271-058-4	68514-79-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-307-00-7	Naphtha (petroleum, full-range reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 35 °C to 230 °C (95 °F to 446 °F).]	272-895-8	68919-37-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-308-00-2	Naphtha (petroleum), catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 30 °C to 220 °C (90 °F to 430 °F). It contains a relatively large proportion of aromatic and branched chain hydrocarbons. This stream may contain 10 vol.% or more benzene.]	273-271-8	68955-35-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-309-00-8	Distillates (petroleum), catalytic reformed hydrotreated light, C_{8-12} arom. fraction; Low boiling point cat-reformed naphtha; [A complex combination of alkylbenzenes obtained by the catalytic reforming of petroleum naphtha. It consists predominantly of alkylbenzenes having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 160 °C to 180 °C (320 °F to 356 °F).]	285-509-8	85116-58-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-310-00-3	Aromatic hydrocarbons, C_8, catalytic reforming-derived; Low boiling point cat-reformed naphtha	295-279-0	91995-18-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-311-00-9	Aromatic hydrocarbons, C_{7-12}, C_8-rich; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} (primarily C_8) and can contain nonaromatic hydrocarbons, both boiling in the range of approximately 130 °C to 200 °C (266 °F to 392 °F).]	297-401-8	93571-75-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-312-00-4	Gasoline, C_{5-11}, high-octane stabilized reformed; Low boiling point cat-reformed naphtha; [A complex high octane combination of hydrocarbons obtained by the catalytic dehydrogenation of a predominantly naphthenic naphtha. It consists predominantly of aromatics and non-aromatics having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 45 °C to 185 °C (113 °F to 365 °F).]	297-458-9	93572-29-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-313-00-X	Hydrocarbons, C_{7-12}, $C_{>9}$-arom.-rich, reforming heavy fraction; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of nonaromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 120 °C to 210 °C (248 °F to 380 °F) and C_9 and higher aromatic hydrocarbons.]	297-465-7	93572-35-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-314-00-5	Hydrocarbons, C_{5-11}, nonaroms.-rich, reforming light fraction; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of nonaromatic hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 35 °C to 125 °C (94 °F to 257 °F), benzene and toluene.]	297-466-2	93572-36-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-315-00-0	Foots oil (petroleum), silicic acid-treated; Foots oil; [A complex combination of hydrocarbons obtained by the treatment of Foots oil with silicic acid for removal of trace constituents and impurities. It consists predominantly of straight chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-127-6	97862-77-6	Carc. 1B	H350 H304	GHS08 Dgr	H350 H304			H L
649-316-00-6	Naphtha (petroleum), light thermal cracked; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons from distillation of products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_8 and boiling in the range of approximately minus 10 °C to 130 °C (14 °F to 266 °F).]	265-075-6	64741-74-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-317-00-1	Naphtha (petroleum), heavy thermal cracked; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons from distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 220 °C (148 °F to 428 °F).]	265-085-0	64741-83-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-318-00-7	Distillates (petroleum), heavy arom.; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons from the distillation of the products from the thermal cracking of ethane and propane. This higher boiling fraction consists predominantly of C_5-C_7 aromatic hydrocarbons with some unsaturated aliphatic hydrocarbons having carbon number predominantly of C_5. This stream may contain benzene.]	267-563-4	67891-79-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-319-00-2	Distillates (petroleum), light arom.; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons from the distillation of the products from the thermal cracking of ethane and propane. This lower boiling fraction consists predominantly of C_5-C_7 aromatic hydrocarbons with some unsaturated aliphatic hydrocarbons having a carbon number predominantly of C_5. This stream may contain benzene.]	267-565-5	67891-80-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-320-00-8	Distillates (petroleum), naphtha-raffinate pyrolyzate-derived, gasoline-blending; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons obtained by the pyrolysis fractionation at 816 °C (1500 °F) of naphtha and raffinate. It consists predominantly of hydrocarbons having a carbon number of C_9 and boiling at approximately 204 °C (400 °F).]	270-344-6	68425-29-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-321-00-3	Aromatic hydrocarbons, C_{6-8}, naphtha-raffinate pyrolyzate-derived; Low boiling point thermally cracked naphtha; A complex combination of hydrocarbons obtained by the fractionation pyrolysis at 816 °C (1500 °F) of naphtha and raffinate. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8, including benzene.]	270-658-3	68475-70-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-322-00-9	Distillates (petroleum), thermal cracked naphtha and gas oil; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by distillation of thermally cracked naphtha and/or gas oil. It consists predominantly of olefinic hydrocarbons having a carbon number of C_5 and boiling in the range of approximately 33 °C to 60 °C (91 °F to 140 °F).]	271-631-9	68603-00-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-323-00-4	Distillates (petroleum), thermal cracked naphtha and gas oil, C_5-dimer-contg.; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the extractive distillation of thermal cracked naphtha and/or gas oil. It consists predominantly of hydrocarbons having a carbon number of C_5 with some dimerized C_5 olefins and boiling in the range of approximately 33 °C to 184 °C (91 °F to 363 °F).]	271-632-4	68603-01-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-324-00-X	Distillates (petroleum), thermal cracked naphtha and gas oil, extractive; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the extractive distillation of thermal cracked naphtha and/or gas oil. It consists of paraffinic and olefinic hydrocarbons, predominantly isoamylenes such as 2-methyl-1-butene and 2-methyl-2-butene and boiling in the range of approximately 31 °C to 40 °C (88 °F to 104 °F).]	271-634-5	68603-03-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-325-00-5	Distillates (petroleum), light thermal cracked, debutanized arom.; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a thermal cracking process. It consists predominantly of aromatic hydrocarbons, primarily benzene.]	273-266-0	68955-29-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-326-00-0	Naphtha (petroleum), light thermal cracked, sweetened; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate from the high temperature thermal cracking of heavy oil fractions to a sweetening process to convert mercaptans. It consists predominantly of aromatics, olefins and saturated hydrocarbons boiling in the range of approximately 20 °C to 100 °C (68 °F to 212 °F).]	295-447-3	92045-65-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-327-00-6	Naphtha (petroleum), hydrotreated heavy; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{13} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-150-3	64742-48-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-328-00-1	Naphtha (petroleum), hydrotreated light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-151-9	64742-49-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-329-00-7	Naphtha (petroleum), hydrodesulfurized light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-178-6	64742-73-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-330-00-2	Naphtha (petroleum), hydrodesulfurized heavy; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-185-4	64742-82-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-331-00-8	Distillates (petroleum), hydrotreated middle, intermediate boiling; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by the distillation of products from a middle distillate hydrotreating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{10} and boiling in the range of approximately 127 °C to 188 °C (262 °F to 370 °F).]	270-092-7	68410-96-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-332-00-3	Distillates (petroleum), light distillate hydrotreating process, low-boiling; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by the distillation of products from the light distillate hydrotreating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_9 and boiling in the range of approximately 3 °C to 194 °C (37 °F to 382 °F).]	270-093-2	68410-97-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-333-00-9	Distillates (petroleum), hydrotreated heavy naphtha, deisohexanizer overheads; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by distillation of the products from a heavy naphtha hydrotreating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 49 °C to 68 °C (- 57 °F to 155 °F).]	270-094-8	68410-98-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-334-00-4	Solvent naphtha (petroleum), light arom., hydrotreated; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F).]	270-988-8	68512-78-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-335-00-X	Naphtha (petroleum), hydrodesulfurized thermal cracked light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by fractionation of hydrodesulfurized thermal cracker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 23 °C to 195 °C (73 °F to 383 °F).]	285-511-9	85116-60-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-336-00-5	Naphtha (petroleum), hydrotreated light, cycloalkane-contg.; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from the distillation of a petroleum fraction. It consists predominantly of alkanes and cycloalkanes boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	285-512-4	85116-61-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-337-00-0	Naphtha (petroleum), heavy steam-cracked, hydrogenated; Low boiling point hydrogen treated naphtha	295-432-1	92045-51-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-338-00-6	Naphtha (petroleum), hydrodesulfurized full-range; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately 30 °C to 250 °C (86 °F to 482 °F).]	295-433-7	92045-52-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-339-00-1	Naphtha (petroleum), hydrotreated light steam-cracked; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction, derived from a pyrolysis process, with hydrogen in the presence of a catalyst. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F).]	295-438-4	92045-57-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-340-00-7	Hydrocarbons, C_{4-12}, naphtha-cracking, hydrotreated; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by distillation from the product of a naphtha steam cracking process and subsequent catalytic selective hydrogenation of gum formers. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 30 °C to 230 °C (86 °F to 446 °F).]	295-443-1	92045-61-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-341-00-2	Solvent naphtha (petroleum), hydrotreated light naphthenic; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists predominantly of cycloparaffinic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_7 and boiling in the range of approximately 73 °C to 85 °C (163 °F to 185 °F).]	295-529-9	92062-15-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-342-00-8	Naphtha (petroleum), light steam-cracked, hydrogenated; Low boiling point hydrogen treated naphtha; [A complex comination of hydrocarbons produced from the separation and subsequent hydrogenation of the products of a steam-cracking process to produce ethylene. It consists predominantly of saturated and unsaturated paraffins, cyclic paraffins and cyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10} and boiling in the range of approximately 50 °C to 200 °C (122 °F to 392 °F). The proportion of benzene hydrocarbons may vary up to 30 wt. % and the stream may also contain small amounts of sulphur and oxygenated compounds.]	296-942-7	93165-55-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-343-00-3	Hydrocarbons, $C_{6\text{-}11}$, hydrotreated, dearomatized; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-852-0	93763-33-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-344-00-9	Hydrocarbons, $C_{9\text{-}12}$, hydrotreated, dearomatized; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-853-6	93763-34-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-345-00-4	Stoddard solvent; Low boiling point naphtha — unspecified; [A colourless, refined petroleum distillate that is free from rancid or objectionable odors and that boils in a range of approximately 300 °F to 400 °F.]	232-489-3	8052-41-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-346-00-X	Natural gas condensates (petroleum); Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated as a liquid from natural gas in a surface separator by retrograde condensation. It consists mainly of hydrocarbons having carbon numbers predominantly in the range of C_2 to C_{20}. It is a liquid at atmospheric temperature and pressure.]	265-047-3	64741-47-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-347-00-5	Natural gas (petroleum), raw liq. mix; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated as a liquid from natural gas in a gas recycling plant by processes such as refrigeration or absorption. It consists mainly of saturated aliphatic hydrocarbons having carbon numbers in the range of C_2 through C_8.]	265-048-9	64741-48-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-348-00-0	Naphtha (petroleum), light hydrocracked; Low boiling naphtha — unspecified; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10}, and boiling in the range of approximately minus 20 °C to 180 °C (- 4 °F to 356 °F).]	265-071-4	64741-69-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-349-00-6	Naphtha (petroleum), heavy hydrocracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12}, and boiling in the range of approximately 65 °C to 230 °C (148 °F to 446 °F).]	265-079-8	64741-78-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-350-00-1	Naphtha (petroleum), sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately minus 10 °C to 230 °C (14 °F to 446 °F).]	265-089-2	64741-87-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-351-00-7	Naphtha (petroleum), acid-treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-115-2	64742-15-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-352-00-2	Naphtha (petroleum), chemically neutralized heavy; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-122-0	64742-22-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-353-00-8	Naphtha (petroleum), chemically neutralized light; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-123-6	64742-23-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-354-00-3	Naphtha (petroleum), catalytic dewaxed; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the catalytic dewaxing of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 35 °C to 230 °C (95 °F to 446 °F).]	265-170-2	64742-66-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-355-00-9	Naphtha (petroleum), light steam-cracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of the products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F). This stream is likely to contain 10 vol.% or more benzene.]	265-187-5	64742-83-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-356-00-4	Solvent naphtha (petroleum), light arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from distillation of aromatic streams. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F).]	265-199-0	64742-95-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-357-00-X	Aromatic hydrocarbons, C_{6-10}, acid-treated, neutralized; Low boiling point naphtha — unspecified	268-618-5	68131-49-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-358-00-5	Distillates (petroleum), C_{3-5}, 2-methyl-2-butene-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons from the distillation of hydrocarbons usually ranging in carbon numbers from C_3 through C_5, predominantly isopentane and 3-methyl-1-butene. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly 2-methyl-2-butene.]	270-725-7	68477-34-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-359-00-0	Distillates (petroleum), polymd. steam-cracked petroleum distillates, C_{5-12} fraction; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the distillation of polymerized steam-cracked petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12}.]]	270-735-1	68477-50-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-360-00-6	Distillates (petroleum), steam-cracked, C_{5-12} fraction; Low boiling point naphtha — unspecified; [A complex combination of organic compounds obtained by the distillation of products from a steam cracking process. It consists of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12}.]	270-736-7	68477-53-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-361-00-1	Distillates (petroleum), steam-cracked, C_{5-10} fraction, mixed with light steam-cracked petroleum naphtha C_5 fraction; Low boiling point naphtha — unspecified	270-738-8	68477-55-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-362-00-7	Extracts (petroleum), cold-acid, C_{4-6}; Low boiling point naphtha — unspecified; [A complex combination of organic compounds produced by cold acid unit extraction of saturated and unsaturated aliphatic hydrocarbons usually ranging in carbon numbers from C_3 through C_6, predominantly pentanes and amylenes. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_4 through C_6, predominantly C_5.]	270-741-4	68477-61-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-363-00-2	Distillates (petroleum), dependentanizer overheads; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from a catalytic cracked gas stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6.]	270-771-8	68477-89-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-364-00-8	Residues (petroleum), butane splitter bottoms; Low boiling point naphtha — unspecified; [A complex residuum from the distillation of butane stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6.]	270-791-7	68478-12-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-365-00-3	Residual oils (petroleum), deisobutanizer tower; Low boiling point naphtha — unspecified; [A complex residuum from the atmospheric distillation of the butane-butylene stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6.]	270-795-9	68478-16-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-366-00-9	Naphtha (petroleum), full-range coker; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by the distillation of products from a fluid coker. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{15} and boiling in the range of approximately 43 °C to 250 °C (110 °F to 500 °F).]	270-991-4	68513-02-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-367-00-4	Naphtha (petroleum), steam-cracked middle arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by the distillation of products from a steam-cracking process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 130 °C to 220 °C (266 °F to 428 °F).]]	271-138-9	68516-20-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-368-00-X	Naphtha (petroleum), clay-treated full-range straight-run; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons resulting from treatment of full-range straight-run naphtha with natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately - 20 °C to 220 °C (- 4 °F to 429 °F).]	271-262-3	68527-21-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-369-00-5	Naphtha (petroleum), clay-treated light straight-run; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons resulting from treatment of light straight-run naphtha with a natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities, present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{10} and boiling in the range of approximately 93 °C to 180 °C (200 °F to 356 °F.]	271-263-9	68527-22-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-370-00-0	Naphtha (petroleum), light steam-cracked arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by distillation of products from a steam-cracking process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_9 and boiling in the range of approximately 110 °C to 165 °C (230 °F to 329 °F).]	271-264-4	68527-23-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-371-00-6	Naphtha (petroleum), light steam-cracked, debenzenized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by distillation of products from a steam-cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 80 °C to 218 °C (176 °F to 424 °F).]	271-266-5	68527-26-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-372-00-1	Naphtha (petroleum), arom.-contg.; Low boiling point naphtha — unspecified	271-635-0	68603-08-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-373-00-7	Gasoline, pyrolysis, debutanizer bottoms; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the fractionation of depropanizer bottoms. It consists of hydrocarbons having carbon numbers predominantly greater than C_5.]	271-726-5	68606-10-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-374-00-2	Naphtha (petroleum), light, sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 20 °C to 100 °C (- 4 °F to 212 °F).]	272-206-0	68783-66-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-375-00-8	Natural gas condensates; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated and/or condensed from natural gas during transportation and collected at the wellhead and/or from the production, gathering, transmission, and distribution pipelines in deeps, scrubbers, etc. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_8.]	272-896-3	68919-39-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H J
649-376-00-3	Distillates (petroleum), naphtha unifiner stripper; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by stripping the products from the naphtha unifiner. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	272-932-8	68921-09-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-377-00-9	Naphtha (petroleum), catalytic reformed light, arom.-free fraction; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons remaining after removal of aromatic compounds from catalytic reformed light naphtha in a selective absorption process. It consists predominantly of paraffinic and cyclic compounds having carbon numbers predominantly in the range of C_5 to C_8 and boiling in the range of approximately 66 °C to 121 °C (151 °F to 250 °F).]	285-510-3	85116-59-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-378-00-4	Gasoline; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons consisting primarily of paraffins, cycloparaffins, aromatic and olefinic hydrocarbons having carbon numbers predominantly greater than C_3 and boiling in the range of 30 °C to 260 °C (86 °F to 500 °F).]	289-220-8	86290-81-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-379-00-X	Aromatic hydrocarbons, C_{7-8}, dealkylation products, distn. residues; Low boiling point naphtha — unspecified	292-698-0	90989-42-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-380-00-5	Hydrocarbons, C_{4-6}, depentanizer lights, arom. hydrotreater; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the depentanizer column before hydrotreatment of the aromatic charges. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6, predominantly pentanes and pentenes, and boiling in the range of approximately 25 °C to 40 °C (77 °F to 104 °F).]	295-298-4	91995-38-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-381-00-0	Distillates (petroleum), heat-soaked steam-cracked naphtha, C_5-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of heat-soaked steam-cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_4 through C_6, predominantly C_5.]	295-302-4	91995-41-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-382-00-6	Extracts (petroleum), catalytic reformed light naphtha solvent; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as the extract from the solvent extraction of a catalytically reformed petroleum cut. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_8 and boiling in the range of approximately 100 °C to 200 °C (212 °F to 392 °F).]	295-331-2	91995-68-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-383-00-1	Naphtha (petroleum), hydrodesulfurized light, dearomatized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of hydrodesulfurized and dearomatized light petroleum fractions. It consists predominantly of C_7 paraffins and cycloparaffins boiling in a range of approximately 90 °C to 100 °C (194 °F to 212 °F).]	295-434-2	92045-53-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-384-00-7	Naphtha (petroleum), light, C_5-rich, sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_5, predominantly C_5, and boiling in the range of approximately minus 10 °C to 35 °C (14 °F to 95 °F).]	295-442-6	92045-60-8	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-385-00-2	Hydrocarbons, C_{8-11}, naphtha-cracking, toluene cut; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation from prehydrogenated cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{11} and boiling in the range of approximately 130 °C to 205 °C (266 °F to 401 °F).]	295-444-7	92045-62-0	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-386-00-8	Hydrocarbons, C_{4-11}, naphtha-cracking, arom.-free; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from prehydrogenated cracked naphtha after distillative separation of benzene- and toluene-containing hydrocarbon cuts and a higher boiling fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range C_4 through C_{11} and boiling in the range of approximately 30 °C to 205 °C (86 °F to 401 °F).]	295-445-2	92045-63-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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649-387-00-3	Naphtha (petroleum), light heat-soaked, steam-cracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the fractionation of steam cracked naphtha after recovery from a heat soaking process. It consists predominantly of hydrocarbons having a carbon numbers predominantly in the range of C_4 through C_6 and boiling in the range of approximately 0 °C to 80 °C (32 °F to 176 °F.]	296-028-8	92201-97-3	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-388-00-9	Distillates (petroleum), C_6-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the distillation of a petroleum feedstock. It consists predominantly of hydrocarbons having carbon numbers of C_5 through C_7, rich in C_6, and boiling in the range of approximately 60 °C to 70 °C (140 °F to 158 °F).]	296-903-4	93165-19-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-389-00-4	Gasoline, pyrolysis, hydrogenated; Low boiling point naphtha — unspecified; [A distillation fraction from the hydrogenation of pyrolysis gasoline boiling in the range of approximately 20 °C to 200 °C (68 °F to 392 °F).]	302-639-3	94114-03-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-390-00-X	Distillates (petroleum), steam-cracked, C_{8-12} fraction, polymd., distn. lights; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of the polymerized C_8 through C_{12} fraction from steam-cracked petroleum distillates. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12}.]	305-750-5	95009-23-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-391-00-5	Extracts (petroleum) heavy naphtha solvent, clay-treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment of heavy naphthic solvent petroleum extract with bleaching earth. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{18} and boiling in the range of approximately 80 °C to 180 °C (175 °F to 356 °F).]	308-261-5	97926-43-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-392-00-0	Naphtha (petroleum), light steam-cracked, debenzenized, thermally treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment and distillation of debenzenized light steam-cracked petroleum naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 95 °C to 200 °C (203 °F to 392 °F).]	308-713-1	98219-46-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-393-00-6	Naphtha (petroleum), light steam-cracked, thermally treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment and distillation of light steam-cracked petroleum naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_6 and boiling in the range of approximately 35 °C to 80 °C (95 °F to 176 °F).]	308-714-7	98219-47-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-394-00-1	Distillates (petroleum), C_{7-9}, C_8-rich, hydrodesulfurized dearomatized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of petroleum light fraction, hydrodesulfurized and dearomatized. It consists predominantly of hydrocarbons having carbon numbers in the range of C_7 through C_9, predominantly C_8 paraffins and cycloparaffins, boiling in the range of approximately 120 °C to 130 °C (248 °F to 266 °F).]	309-862-5	101316-56-7	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-395-00-7	Hydrocarbons, C_{6-8}, hydrogenated sorption-dearomatized, toluene raffination; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained during the sorptions of toluene from a hydrocarbon fraction from cracked gasoline treated with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8 and boiling in the range of approximately 80 °C to 135 °C (176 °F to 275 °F).]	309-870-9	101316-66-9	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-396-00-2	Naphtha (petroleum), hydrodesulfurized full-range coker; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized coker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 23 °C to 196 °C (73 °F to 385 °F).]	309-879-8	101316-76-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-397-00-8	Naphtha (petroleum), sweetened light; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_8 and boiling in the range of approximately 20 °C to 130 °C (68 °F to 266 °F).]	309-976-5	101795-01-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-398-00-3	Hydrocarbons, C_{3-6}, C_5-rich, steam-cracked naphtha; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of steam-cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly C_5.]	310-012-0	102110-14-5	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-399-00-9	Hydrocarbons, C_5-rich, dicyclopentadiene-contg.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of the products from a stream-cracking process. It consists predominantly of hydrocarbons having carbon numbers of C_5 and dicyclopentadiene and boiling in the range of approximately 30 °C to 170 °C (86 °F to 338 °F).]	310-013-6	102110-15-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-400-00-2	Residues (petroleum), steam-cracked light, arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of the products of steam cracking or similar processes after taking off the very light products resulting in a residue starting with hydrocarbons having carbon numbers greater than C_5. It consists predominantly of aromatic hydrocarbons having carbon numbers greater than C_5 and boiling above approximately 40 °C (104 °F).]	310-057-6	102110-55-4	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-401-00-8	Hydrocarbons, $C_{\geq 5}$, C_{5-6}-rich; Low boiling point naphtha — unspecified	270-690-8	68476-50-6	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-402-00-3	Hydrocarbons, C_5-rich; Low boiling point naphtha — unspecified	270-695-5	68476-55-1	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-403-00-9	Aromatic hydrocarbons, C_{8-10}; Low boiling point naphtha — unspecified	292-695-4	90989-39-2	Carc. 1B Asp. Tox. 1	H350 H304	GHS08 Dgr	H350 H304			H P
649-404-00-4	Kerosine (petroleum); Straight run kerosine; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (320 °F to 554 °F).]	232-366-4	8008-20-6	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-405-00-X	Solvent naphtha (petroleum), medium aliph.; Straight run kerosine; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{12} and boiling in the range of approximately 140 °C to 220 °C (284 °F to 428 °F).]	265-191-7	64742-88-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-406-00-5	Solvent naphtha (petroleum) heavy aliph.; Straight run kerosine; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{16} and boiling in the range of approximately 190 °C to 290 °C (374 °F to 554 °F).]	265-200-4	64742-96-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-407-00-0	Kerosine (petroleum), straight-run wide-cut; Straight run kerosine; [A complex combination of hydrocarbons obtained as a wide cut hydrocarbon fuel cut from atmospheric distillation and boiling in the range of approximately 70 °C to 220 °C (158 °F to 428 °F).]	295-418-5	92045-37-9	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-408-00-6	Distillates (petroleum), steam-cracked; Cracked kerosine; [A complex combination of hydrocarbons obtained by the distillation of the products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{16} and boiling in the range of approximately 90 °C to 290 °C (190 °F to 554 °F).]	265-194-3	64742-91-2	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-409-00-1	Distillates (petroleum), cracked stripped steam-cracked petroleum distillates, C_{8-10} fraction; Cracked kerosine; [A complex combination of hydrocarbons obtained by distilling cracked stripped steam-cracked distillates. It consists of hydro-carbons having carbon numbers in the range of C_8 through C_{10} and boiling in the range of approximately 129 °C to 194 °C (264 °F to 382 °F).]	270-728-3	68477-39-4	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-410-00-7	Distillates (petroleum), cracked stripped steam-cracked petroleum distillates, C_{10-12} fraction; Cracked kerosine; [A complex combination of hydrocarbons obtained by distilling cracked stripped steam-cracked distillates. It consists predominantly of aromatic hydrocarbons having carbon numbers in the range of C_{10} through C_{12}.]	270-729-9	68477-40-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-411-00-2	Distillates (petroleum), steam-cracked, C_{8-12} fraction; Cracked kerosine; [A complex combination of organic compounds obtained by the distillation of products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12}.]	270-737-2	68477-54-3	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-412-00-8	Kerosine (petroleum), hydrodesulfurized thermal cracked; Cracked kerosine; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized thermal cracker distillate. It consists predominantly of hydrocarbons predominantly in the range of C_8 to C_{16} and boiling in the range of approximately 120 °C to 283 °C (284 °F to 541 °F).]	285-507-7	85116-55-8	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-413-00-3	Aromtic hydrocarbons, $C_{\geq 10}$, steam-cracking, hydrotreated; Cracked kerosine; [A complex combination of hydrocarbons produced by the distillation of the products from a steam cracking process treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly greater than C_{10} and boiling in the range of approximately 150 °C to 320 °C (302 °F to 608 °F).]	292-621-0	90640-98-5	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-414-00-9	Naphtha (petroleum), steam-cracked, hydrotreated, C_{9-10}-arom.-rich; Cracked kerosine; [A complex combination of hydrocarbons produced by the distillation of the products from a steam cracking process thereafter treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers in the range of C_9 through C_{10} and boiling in the range of approximately 140 °C to 200 °C (284 °F to 392 °F).]	292-637-8	90641-13-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-415-00-4	Distillates (petroleum), thermal-cracked, alkylarom. hydrocarbon-rich; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of thermal-cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F.]	309-866-7	101316-61-4	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-416-00-X	Distillates (petroleum), catalytic cracked heavy tar light; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of catalytic cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F).]	309-938-8	101631-13-4	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-417-00-5	Solvent naphtha (petroleum), hydrocracked heavy arom.; Cracked kerosine; [A complex combination of hydrocarbons obtained by the distillation of hydrocracked petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 235 °C to 290 °C (455 °F to 554 °F).]	309-881-9	101316-80-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-418-00-0	Distillates (petroleum), steam-cracked heavy tar light; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of steam cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F).]	309-940-9	101631-15-6	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-419-00-6	Distillates (petroleum), alkylate; Kerosine — unspecified; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_{11} through C_{17} and boiling in the range of approximately 205 °C to 320 °C (401 °F to 608 °F).]	265-074-0	64741-73-7	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-420-00-1	Extracts (petroleum), heavy naphtha solvent; Kerosine — unspecified; [A complex combination of hydrocarbons obtained as the extract from a solvent extraction process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 220 °C (194 °F to 428 °F).]	265-099-7	64741-98-6	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-421-00-7	Distillates (petroleum), chemically neutralized light; Kerosine — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-132-5	64742-31-0	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-422-00-2	Distillates (petroleum), hydrotreated light; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-149-8	64742-47-8	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-423-00-8	Kerosine (petroleum), hydrodesulfurized; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-184-9	64742-81-0	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-424-00-3	Solvent naphtha (petroleum), heavy arom.; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from distillation of aromatic streams. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 165 °C to 290 °C (330 °F to 554 °F).]	265-198-5	64742-94-5	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-425-00-9	Naphtha (petroleum), heavy coker; Kerosine — unspecified; [A complex combination of hydrocarbons from the distillation of products from a fluid coker. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{15} and boiling in the range of approximately 157 °C to 288 °C (315 °F to 550 °F).]	269-778-9	68333-23-3	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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649-426-00-4	Naphtha (petroleum), catalytic reformed hydrodesulfurized heavy, arom. fraction; Kerosine — unspecified; [A complex combination of hydrocarbons produced by fractionation from catalytically reformed hydrodesulfurized naphtha. It consists predominantly of aromatic hydrocarbons having carbon numbers predominently in the range of C_7 to C_{13} and boiling in the range of approximately 98 °C to 218 °C (208 °F to 424 °F).]	285-508-2	85116-57-0	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-427-00-X	Kerosine (petroleum), sweetened; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of 130 °C to 290 °C (266 °F to 554 °F).]	294-799-5	91770-15-9	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-428-00-5	Kerosine (petroleum), solvent-refined sweetened; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by solvent refining and sweetening and boiling in the range of approximately 150 °C to 260 °C (302 °F to 500 °F).]	295-416-4	92045-36-8	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-429-00-0	Hydrocarbons, C_{9-16}, hydrotreated, dearomatized; Kerosine — unspecified; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-854-1	93763-35-0	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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649-430-00-6	Kerosine (petroleum), solvent-refined hydrodesulfurized; Kerosine — unspecified	307-033-2	97488-94-3	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-431-00-1	Distillates (petroleum), hydrodesulfurized full-range middle coker; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized coker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{16} and boiling in the range of approximately 120 °C to 283 °C (248 °F to 541 °F).]	309-864-6	101316-58-9	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-432-00-7	Solvent naphtha (petroleum), hydrodesulfurized heavy arom.; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by the catalytic hydrodesulfurization of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{13} and boiling in the range of approximately 180 °C to 240 °C (356 °F to 464 °F).]	309-882-4	101316-81-8	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-433-00-2	Solvent naphtha (petroleum), hydrodesulfurized medium; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by the catalytic hydrodesulfurization of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{13} and boiling in the range of approximately 175 °C to 220 °C (347 °F to 428 °F).]	309-884-5	101316-82-9	Asp. Tox. 1	H304	GHS08 Dgr	H304			H

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649-434-00-8	Kerosine (petroleum), hydrotreated; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from the distillation of petroleum and subsequent hydrotreatment. It consists predominantly of alkanes, cycloalkanes and alkylbenzenes having carbon numbers predominantly in the range of C_{12} through C_{16} and boiling in the range of approximately 230 °C to 270 °C (446 °F to 518 °F).]	309-944-0	101631-19-0	Asp. Tox. 1	H304	GHS08 Dgr	H304			H
649-435-00-3	Distillates (petroleum), light catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F). It contains a relatively large proportion of bicyclic aromatic hydrocarbons.]	265-060-4	64741-59-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-436-00-9	Distillates (petroleum), intermediate catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{30} and boiling in the range of approximately 205 °C to 450 °C (401 °F to 842 °F). It contains a relatively large proportion of tricyclic aromatic hydrocarbons.]	265-062-5	64741-60-2	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-437-00-4	Distillates (petroleum), light hydrocracked; Cracked gasoil; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{18} and boiling in the range of approximately 160 °C to 320 °C (320 °F to 608 °F).]	265-078-2	64741-77-1	Carc. 2	H351	GHS08 Wng	H351			H
649-438-00-X	Distillates (petroleum), light thermal cracked; Cracked gasoil; [A complex combination of hydrocarbons from the distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{22} and boiling in the range of approximately 160 °C to 370 °C (320 °F to 698 °F).]	265-084-5	64741-82-8	Carc. 1B	H350	GHS08 Dgr	H350			H
649-439-00-5	Distillates (petroleum), hydrodesulfurized light catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons obtained by treating light catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F). It contains a relatively large proportion of bicyclic aromatic hydrocarbons.]	269-781-5	68333-25-5	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-440-00-0	Distillates (petroleum), light steam-cracked naphtha; Cracked gasoil; [A complex combination of hydrocarbons from the multiple distillation of products from a steam cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{18}.]	270-662-5	68475-80-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-441-00-6	Distillates (petroleum), cracked steam-cracked petroleum distillates; Cracked gasoil; [A complex combination of hydrocarbons produced by distilling cracked steam cracked distillate and/or its fractionation products. It consists of hydrocarbons having carbon numbers predominently in the range of C_{10} to low molecular weight polymers.]	270-727-8	68477-38-3	Carc. 1B	H350	GHS08 Dgr	H350			H
649-442-00-1	Gas oils (petroleum), steam-cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by distillation of the products from a steam cracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_9 and boiling in the range of from approximately 205 °C to 400 °C (400 °F to 752 °F).]	271-260-2	68527-18-4	Carc. 1B	H350	GHS08 Dgr	H350			H
649-443-00-7	Distillates (petroleum), hydrodesulfurized thermal cracked middle; Cracked gasoil; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized themal cracker distillate stocks. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{11} to C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	285-505-6	85116-53-6	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-444-00-2	Gas oils (petroleum), thermal-cracked, hydrodesulfurized; Cracked gasoil	295-411-7	92045-29-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-445-00-8	Residues (petroleum), hydrogenated steam-cracked naphtha; Cracked gasoil; [A complex combination of hydrocarbons obtained as a residual fraction from the distillation of hydrotreated steam-cracked naphtha. It consists predominantly of hydrocarbons boiling in the range of approximately 200 °C to 350 °C (32 °F to 662 °F).]	295-514-7	92062-00-5	Carc. 1B	H350	GHS08 Dgr	H350			H
649-446-00-3	Residues (petroleum), steam-cracked naphtha distn.; Cracked gasoil; [A complex combination of hydrocarbons obtained as a column bottom from the separation of effluents from steam cracking naphtha at a high temperature. It boils in the range of approximately 147 °C to 300 °C (297 °F to 572 °F) and produces a finished oil having a viscosity of 18cSt at 50 °C.]	295-517-3	92062-04-9	Carc. 1B	H350	GHS08 Dgr	H350			H
649-447-00-9	Distillates (petroleum), light catalytic cracked, thermally degraded; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process which has been used as a heat transfer fluid. It consists predominantly of hydrocarbons boiling in the range of approximately 190 °C to 340 °C (374 °F to 644 °F). This stream is likely to contain organic sulfur compounds.]	295-991-1	92201-60-0	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-448-00-4	Residues (petroleum), steam-cracked heat-soaked naphtha; Cracked gasoil; [A complex combination of hydrocarbons obtained as residue from the distillation of steam cracked heat soaked naphtha and boiling in the range of approximately 150 °C to 350 °C (302 °F to 662 °F).]	297-905-8	93763-85-0	Carc. 1B	H350	GHS08 Dgr	H350			H
649-449-00-X	Hydrocarbons, C_{16-20}, solvent-dewaxed hydrocracked paraffinic distn. residue; Cracked gasoil; [A complex combination of hydrocarbons obtained by solvent dewaxing of a distillation residue from a hydrocracked paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{20} and boiling in the range of approximately 360 °C to 500 °C (680 °F to 932 °F). It produces a finished oil having a viscosity of 4,5 cSt at approximately 100 °C (212 °F).]	307-662-2	97675-88-2	Carc. 2	H351	GHS08 Wng	H351			H
649-450-00-5	Gas oils (petroleum), light vacuum, thermal-cracked hydrodesulfurized; Cracked gasoil; [A complex combination of hydrocarbons obtained by catalytic dehydrosulfurization of thermal-cracked light vacuum petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{20} and boiling in the range of approximately 270 °C to 370 °C (518 °F to 698 °F).]	308-278-8	97926-59-5	Carc. 1B	H350	GHS08 Dgr	H350			H

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-451-00-0	Distillates (petroleum), hydrodesulfurized middle coker; Cracked gasoil; [A complex combination of hydrocarbons by fractionation from hydrodesulfurised coker distillate stocks. Is consists of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{21} and boiling in the range of approximately 200 °C to 360 °C (392 °F to 680 °F).]	309-865-1	101316-59-0	Carc. 1B	H350	GHS08 Dgr	H350			H
649-452-00-6	Distillates (petroleum), heavy steam-cracked; Cracked gasoil; [A complex combination of hydrocarbons obtained by distillation of steam cracking heavy residues. It consists predominantly of highly alkylated heavy aromatic hydrocarbons boiling in the range of approximately 250 °C to 400 °C (482 °F to 752 °F).]	309-939-3	101631-14-5	Carc. 1B	H350	GHS08 Dgr	H350			H
649-453-00-1	Distillates (petroleum), heavy hydrocracked; Baseoil — unspecified; [A complex combination of hydrocarbons from the distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers in the range of C_{15}-C_{39} and boiling in the range of approximately 260 °C to 600 °C (500 °F to 1112 °F).]	265-077-7	64741-76-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-454-00-7	Distillates (petroleum), solvent-refined heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C).]	265-090-8	64741-88-4	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-455-00-2	Distillates (petroleum), solvent-refined light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-091-3	64741-89-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-456-00-8	Residual oils (petroleum), solvent deasphalted; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the solvent soluble fraction from C_3-C_4 solvent deasphalting of a residuum. It consists of hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-096-0	64741-95-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-457-00-3	Distillates (petroleum), solvent-refined heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt a 40 °C). It contains relatively few normal paraffins.]	265-097-6	64741-96-4	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-458-00-9	Distillates (petroleum), solvent-refined light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-098-1	64741-97-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-459-00-4	Residual oils (petroleum,) solvent-refined; Baseoil — unspecified; [A complex combination by hydrocarbons obtained as the solvent insoluble fraction from solvent refining of a residuum using a polar organic solvent such as phenol or furfural. It consists of hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-101-6	64742-01-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-460-00-X	Distillates (petroleum), clay-treated paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-137-2	64742-36-5	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-461-00-5	Distillates (petroleum), clay-treated light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-138-8	64742-37-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-462-00-0	Residual oils (petroleum), clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treatment of a residual oil with a natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-143-5	64742-41-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-463-00-6	Distillates (petroleum), clay-treated heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-146-1	64742-44-5	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-464-00-1	Distillates (petroleum), clay-treated light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-147-7	64742-45-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-465-00-7	Distillates (petroleum), hydrotreated heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-155-0	64742-52-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-466-00-2	Distillates (petroleum), hydrotreated light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-156-6	64742-53-6	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-467-00-8	Distillates (petroleum), hydrotreated heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-157-1	64742-54-7	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-468-00-3	Distillates (petroleum), hydrotreated light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-158-7	64742-55-8	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-469-00-9	Distillates (petroleum), solvent-dewaxed light paraffinic; Baseoil — unspecified; [A complex comination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-159-2	64742-56-9	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-470-00-4	Residual oils (petroleum), hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly greater than C_{25} and boiling above approximately 400 °C (752 °F).]	265-160-8	64742-57-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-471-00-X	Residual oils (petroleum), solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of long, branched chain hydrocarbons from a residual oil by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly greater than C_{25} and boiling above approximately 400 °C (752 °F).]	265-166-0	64742-62-7	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-472-00-5	Distillates (petroleum), solvent-dewaxed heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} . through C_{50} and produces a finished oil of not less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-167-6	64742-63-8	Carc. 1B	H350	GHS08 Dgr	H350			H L

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649-473-00-0	Distillates (petroleum), solvent-dewaxed light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly in the range C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-168-1	64742-64-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-474-00-6	Distillates (petroleum), solvent-dewaxed heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity not less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-169-7	64742-65-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-475-00-1	Naphthenic oils (petroleum), catalytic dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-172-3	64742-68-3	Carc. 1B	H350	GHS08 Dgr	H350			H L

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649-476-00-7	Naphthenic oils (petroleum), catalytic dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-173-9	64742-69-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-477-00-2	Paraffin oils (petroleum), catalytic dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C).]	265-174-4	64742-70-7	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-478-00-8	Paraffin oils (petroleum), catalytic dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-176-5	64742-71-8	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-479-00-3	Naphthenic oils (petroleum), complex dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removing straight chain paraffin hydrocarbons as a solid by treatment with an agent such as urea. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-179-1	64742-75-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-480-00-9	Naphthenic oils (petroleum), complex dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-180-7	64742-76-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-481-00-4	Lubricating oils (petroleum), C_{20-50}, hydrotreated neutral oil-based, high-viscosity; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil, heavy vacuum gas oil, and solvent deasphalted residual oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of approximately 112cSt at 40 °C. It contains a relatively large proportion of saturated hydrocarbons.]	276-736-3	72623-85-9	Carc. 1B	H350	GHS08 Dgr	H350			H L

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649-482-00-X	Lubricating oils (petroleum), C_{15-30}, hydrotreated neutral oil-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil and heavy vacuum gas oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity of approximately 15cSt at 40 °C. It contains a relatively large proportion of saturated hydrocabons.]	276-737-9	72623-86-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-483-00-5	Lubricating oils (petroleum), C_{20-50}, hydrotreated neutral oil-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil, heavy vacuum gas oil and solvent deasphalted residual oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of approximately 32cSt at 40 °C. It contains a relatively large proportion of saturated hydrocarbons.]	276-738-4	72623-87-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-484-00-0	Lubricating oils; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from solvent extraction and dewaxing processes. It consists predominantly of saturated hydrocarbons having carbon numbers in the range C_{15} through C_{50}.]	278-012-2	74869-22-0	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-485-00-6	Distillates (petroleum), complex dewaxed heavy paraffinci; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by dewaxing heavy paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of equal to or greater than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	292-613-7	90640-91-8	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-486-00-1	Distillates (petroleum), complex dewaxed light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by dewaxing light paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	292-614-2	90640-92-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-487-00-7	Distillates (petroleum), solvent dewaxed heavy paraffinic, clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating dewaxed heavy paraffinic distillate with neutral or modified clay in either a contacting or percolation process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	292-616-3	90640-94-1	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-488-00-2	Hydrocarbons, C_{20-50}, solvent dewaxed heavy paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons produced by treating dewaxed heavy paraffinic distillate with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	292-617-9	90640-95-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-489-00-8	Distillates (petroleum), solvent dewaxed light paraffinic, clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of dewaxed light paraffinic distillate with natural or modified clay in either a contacting or percolation process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}.]	292-618-4	90640-96-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-490-00-3	Distillates (petroleum), solvent dewaxed light paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons produced by treating a dewaxed light paraffinic distillate with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}.]	292-620-5	90640-97-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-491-00-9	Residual oils (petroleum), hydrotreated solvent dewaxed; Baseoil — unspecified	292-656-1	90669-74-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-492-00-4	Residual oils (petroleum), catalytic dewaxed; Baseoil — unspecified	294-843-3	91770-57-9	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-493-00-X	Distillates (petroleum), dewaxed heavy paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from an intensive treatment of dewaxed distillate by hydrogenation in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{25} through C_{39} and produces a finished oil with a viscosity of approximately 44 cSt at 50 °C.]	295-300-3	91995-39-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-494-00-5	Distillates (petroleum), dewaxed light paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from an intensive treatment of dewaxed distillate by hydrogenation in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{21} through C_{29} and produces a finished oil with a viscosity of approximately 13 cSt at 50 °C.]	295-301-9	91995-40-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-495-00-0	Distillates (petroleum), hydrocracked solvent-refined, dewaxed; Baseoil — unspecified; [A complex combination of liquid hydrocarbons obtained by recrystallization of dewaxed hydrocracked solvent-refined petroleum distillates.]	295-306-6	91995-45-8	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-496-00-6	Distillates (petroleum), solvent-refined light naphthenic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst and removing the aromatic hydrocarbons by solvent extraction. It consists predominantly of naphthenic hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of between 13-15cSt at 40 °C.]	295-316-0	91995-54-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-497-00-1	Lubricating oils (petroleum), C_{17-35}, solvent-extd., dewaxed, hydrotreated; Baseoil — unspecified	295-423-2	92045-42-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-498-00-7	Lubricating oils (petroleum), hydrocracked nonarom. solvent-deparaffined; Baseoil — unspecified	295-424-8	92045-43-7	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-499-00-2	Residual oils (petroleum), hydrocracked acid-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons produced by solvent removal of paraffins from the residue of the distillation of acid-treated, hydrocracked heavy paraffins and boiling approximately above 380 °C (716 °F).]	295-499-7	92061-86-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-500-00-6	Paraffin oils (petroleum), solvent-refined dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from sulfur-containing paraffinic crude oil. It consists predominantly of a solvent refined deparaffinated lubricating oil with a viscosity of 65cSt at 50 °C.]	295-810-6	92129-09-4	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-501-00-1	Lubricating oils (petroleum), base oils, paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by refining of crude oil. It consists predominantly of aromatics, naphthenics and paraffinics and produces a finished oil with a viscosity of 120 SUS at 100 °F (23cSt at 40 °C).]	297-474-6	93572-43-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-502-00-7	Hydrocarbons, hydrocracked paraffinic distn. residues, solvent-dewaxed; Baseoil — unspecified	297-857-8	93763-38-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-503-00-2	Hydrocarbons, C_{20-50}, residual oil hydrogenation vacuum distillate; Baseoil — unspecified	300-257-1	93924-61-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-504-00-8	Distillates (petroleum), solvent-refined hydrotreated heavy; hydrogenated; Baseoil — unspecified	305-588-5	94733-08-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-505-00-3	Distillates (petroleum), solvent-refined hydrocracked light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent dearomatization of the residue of hydrocracked petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{27} and boiling in the range of approximately 370 °C to 450 °C (698 °F to 842 °F).]	305-589-0	94733-09-2	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-506-00-9	Lubricating oils (petroleum), C_{18-40}, solvent-dewaxed hydrocracked distillate-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent deparaffination of the distillation residue from hydrocracked petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{40} and boiling in the range of approximately 370 °C to 550 °C (698 °F to 1022 °F).]	305-594-8	94733-15-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-507-00-4	Lubricating oils (petroleum), C_{18-40}, solvent-dewaxed hydrogenated raffinate-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent deparaffination of the hydrogenated raffinate obtained by solvent extraction of a hydrotreated petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{40} and boiling in the range of approximately 370 °C to 550 °C (698 °F to 1022 °F).]	305-595-3	94733-16-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-508-00-X	Hydrocarbons, C_{13-30}, arom.-rich, solvent-extd. naphthenic distillate; Baseoil — unspecified	305-971-7	95371-04-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-509-00-5	Hydrocarbons, C_{16-32}, arom. rich, solvent-extd. naphthenic distillate; Baseoil — unspecified	305-972-2	95371-05-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-510-00-0	Hydrocarbons, C_{37-68}, dewaxed deasphalted hydrotreated vacuum distn. residues; Baseoil — unspecified	305-974-3	95371-07-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-511-00-6	Hydrocarbons, C_{37-65}, hydrotreated deasphalted vacuum distn. residues; Baseoil — unspecified	305-975-9	95371-08-7	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-512-00-1	Distillates (petroleum), hydrocracked solvent-refined light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by the solvent treatment of a distillate from hydrocracked petroleum distillates. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{27} and boiling in the range of approximately 370 °C to 450 °C (698 °F to 842 °F.]	307-010-7	97488-73-8	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-513-00-7	Distillates (petroleum), solvent-refined hydrogenated heavy; Baseoil — unspecified; [A complex combination of hydrocarbons, obtained by the treatment of a hydrogenated petroleum distillate with a solvent. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{19} through C_{40} and boiling in the range of approximately 390 °C to 550 °C (734 °F to 1022 °F).]	307-011-2	97488-74-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-514-00-2	Lubricating oils (petroleum), C_{18-27}, hydrocracked solvent-dewaxed; Baseoil — unspecified	307-034-8	97488-95-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-515-00-8	Hydrocarbons, C_{17-30}, hydrotreated solvent-deasphalted atm. distn. residue, distn. lights; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of a solvent deasphalted short residue with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{30} and boiling in the range of approximately 300 °C to 400 °C (572 °F to 752 °F). It produces a finished oil having a viscosity of 4cSt at approximately 100 °C (212 °F).]	307-661-7	97675-87-1	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-516-00-3	Hydrocarbons, C_{17-40}, hydrotreated solvent-deasphalted distn. residue, vacuum distn. lights; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the catalytic hydrotreatment of a solvent deasphalted short residue having a viscosity of 8cSt at approximately 100 °C (212 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{40} and boiling in the range of approximately 300 °C to 500 °C (592 °F to 932 °F).]	307-755-8	97722-06-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-517-00-9	Hydrocarbons, C_{13-27}, solvent-extd. light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a viscosity of 9.5cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{27} and boiling in the range of approximately 240 °C to 400 °C (464 °F to 752 °F).]	307-758-4	97722-09-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-518-00-4	Hydrocarbons, C_{14-29}, solvent-extd. light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a viscosity of 16cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{29} and boiling in the range of approximately 250 °C to 425 °C (482 °F to 797 °F).]	307-760-5	97722-10-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-519-00-X	Hydrocarbons, C_{27-42}, dearomatized; Baseoil — unspecified	308-131-8	97862-81-2	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-520-00-5	Hydrocarbons, C_{17-30}, hydrotreated distillates, distn. lights; Baseoil — unspecified	308-132-3	97862-82-3	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-521-00-0	Hydrocarbons, C_{27-45}, naphthenic vacuum distn.; Baseoil — unspecified	308-133-9	97862-83-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-522-00-6	Hydrocarbons, C_{27-45}, dearomatized; Baseoil — unspecified	308-287-7	97926-68-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-523-00-1	Hydrocarbons, C_{20-58}, hydrotreated; Baseoil — unspecified	308-289-8	97926-70-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-524-00-7	Hydrocarbons, C_{27-42}, naphthenic; Baseoil — unspecified	308-290-3	97926-71-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-525-00-2	Residual oils (petroleum), carbon-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of solvent-dewaxed petroleum residual oils with activated charcoal for the removal of trace polar constituents and impurities.]	309-710-8	100684-37-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-526-00-8	Residual oils (petroleum), clay-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treatment of solvent-dewaxed petroleum residual oils with bleaching earth for the removal of trace polar constituents and impurities.]	309-711-3	100684-38-6	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-527-00-3	Lubricating oils (petroleum), $C_{>25}$, solvent-extd., deasphalted, dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of vacuum distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{25} and produces a finished oil with a viscosity in the order of 32cSt to 37cSt at 100 °C (212 °F).]	309-874-0	101316-69-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-528-00-9	Lubricating oils (petroleum), C_{17-32}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{32} and produced a finished oil with a viscosity in the order of 17cSt to 23cSt at 40 °C (104 °F.]	309-875-6	101316-70-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-529-00-4	Lubricating oils (petroleum), C_{20-35}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{35} and produces a finished oil with a viscosity in the order of 37cSt to 44cSt at 40 °C (104 °F).]	309-876-1	101316-71-6	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-530-00-X	Lubricating oils (petroleum), C_{24-50}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{24} through C_{50} and produces a finished oil with a viscosity in the order of 16cSt to 75cSt at 40 °C (104 °F).]	309-877-7	101316-72-7	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-531-00-5	Extracts (petroleum), heavy naphthenic distillate solvent, arom. conc.; Distillate aromatic extract (treated); [An aromatic concentrate produced by adding water to heavy naphthenic distillate solvent extract and extraction solvent.]	272-175-3	68783-00-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-532-00-0	Extracts (petroleum), solvent-refined heavy paraffinic distillate solvent; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from the re-extraction of solvent-refined heavy paraffinic distillate. It consists of saturated and aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	272-180-0	68783-04-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-533-00-6	Extracts (petroleum), heavy paraffinic distillates, solvent-deasphalted; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from a solvent extraction of heavy paraffinic distillate.]	272-342-0	68814-89-1	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-534-00-1	Extracts (petroleum), heavy naphthenic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by treating a heavy naphthenic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 19cSt at 40 °C (100 SUS at 100 °F).]	292-631-5	90641-07-9	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-535-00-7	Extracts (petroleum), heavy paraffinic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons produced by treating a heavy paraffinic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{21} through C_{33} and boiling in the range of approximately 350 °C to 480 °C (662 °F to 896 °F).	292-632-0	90641-08-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-536-00-2	Extracts (petroleum), light paraffinic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons produced by treating a light paraffinic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{26} and boiling in the range of approximately 280 °C to 400 °C (536 °F to 752 °F).]	292-633-6	90641-09-1	Carc. 1B	H350	GHS08 Dgr	H350			H L

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649-537-00-8	Extracts (petroleum), hydrotreated light paraffinic distillate solvent; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from solvent extraction of intermediate paraffinic top solvent distillate that is treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	295-335-4	91995-73-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-538-00-3	Extracts (petroleum), light naphthenic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by treating the extract, obtained from a solvent extraction process, with hydrogen in the presence of a catalyst under conditions primarily to remove sulfur compounds. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}. This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	295-338-0	91995-75-4	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-539-00-9	Extracts (petroleum), light paraffinic distillate solvent, acid-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction of the distillation of an extract from the solvent extraction of light paraffinic top petroleum distillates that is subjected to a sulfuric acid refining. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	295-339-6	91995-76-5	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-540-00-4	Extracts (petroleum), light paraffinic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of a light paraffin distillate and treated with hydrogen to convert the organic sulfur to hydrogen sulfide which is eliminated. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{40} and produces a finished oil with a viscosity of greater than 10cSt at 40 °C.]	295-340-1	91995-77-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-541-00-X	Extracts (petroleum), light vacuum gas oil solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons, obtained by solvent extraction from light vacuum petroleum gas oils and treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	295-342-2	91995-79-8	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-542-00-5	Extracts (petroleum), heavy paraffinic distillate solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contact or percolation process to remove the trace amounts of polar compounds and impurities present. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}. This stream is likely to contain 5 wt.% or more 4-6 membered ring aromatic hydrocarbons.]	296-437-1	92704-08-0	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-543-00-0	Extracts (petroleum), heavy naphthenic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{50} and produces a finished oil with a viscosity of greater than 19cSt at 40 °C.]	297-827-4	93763-10-1	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-544-00-6	Extracts (petroleum), solvent-dewaxed heavy paraffinic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained from a solvent dewaxed petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{50} and produces a finished oil with a viscosity of greater than 19cSt at 40 °C.]	297-829-5	93763-11-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-545-00-1	Extracts (petroleum), light paraffinic distillate solvent, carbon-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction from distillation of an extract recovered by solvent extraction of light paraffinic top petroleum distillate treated with activated charcoal to remove traces of polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	309-672-2	100684-02-4	Carc. 1B	H350	GHS08 Dgr	H350			H L

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				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-546-00-7	Extracts (petroleum), light paraffinic distillate solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction from distillation of an extract recovered by solvent extraction of light paraffinic top petroleum distillates treated with bleaching earth to remove traces of polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	309-673-8	100684-03-5	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-547-00-2	Extracts (petroleum), light vacuum, gas oil solvent, carbon-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of light vacuum petroleum gas oil treated with activated charcoal for the removal of trace polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	309-674-3	100684-04-6	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-548-00-8	Extracts (petroleum), light vacuum gas oil solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of light vacuum petroleum gas oils treated with bleaching earth for removal of trace polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	309-675-9	100684-05-7	Carc. 1B	H350	GHS08 Dgr	H350			H L

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
649-549-00-3	Foots oil (petroleum); Foots oil; [A complex combination of hydrocarbons obtained as the oil fraction from a solvent deoiling or a wax sweating process. It consists predominantly of branched chain hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	265-171-8	64742-67-2	Carc. 1B	H350	GHS08 Dgr	H350			H L
649-550-00-9	Foots oil (petroleum), hydrotreated; Foots oil	295-394-6	92045-12-0	Carc. 1B	H350	GHS08 Dgr	H350			H L
650-002-00-6	turpentine, oil	232-350-7	8006-64-2	Flam. Liq. 3 Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Asp. Tox. 1 Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Chronic 2	H226 H332 H312 H302 H304 H319 H315 H317 H411	GHS02 GHS08 GHS07 GHS09 Dgr	H226 H332 H312 H302 H304 H319 H315 H317 H411			
650-003-00-1	fenson (ISO); 4-chlorophenyl benzenesulphonate;	201-274-6	80-38-6	Acute Tox. 4 * Eye Irrit. 2 Aquatic Chronic 2	H302 H319 H411	GHS07 GHS09 Wng	H302 H319 H411			
650-004-00-7	norbormide (ISO); 5-(α-hydroxy-α-2-pyridylbenzyl)-7-(α-2-pyridylbenzylidene)bicyclo [2.2.1] hept-5-ene-2,3-dicarboximide	213-589-6	991-42-4	Acute Tox. 4 *	H302	GHS07 Wng	H302			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
650-005-00-2	(2*R*,6a*S*,12a*S*)-1,2,6,6a,12,12a-hexahydro-2-isopropenyl-8,9-dimethoxychromeno[3,4-b]furo[2,3-h]chromen-6-one, rotenone	201-501-9	83-79-4	Acute Tox. 3 * Eye Irrit. 2 STOT SE 3 Skin Irrit. 2 Aquatic Acute 1 Aquatic Chronic 1	H301 H319 H335 H315 H400 H410	GHS06 GHS09 Dgr	H301 H319 H335 H315 H410			
650-006-00-8	benquinox (ISO); *p*-benzoquinone 1-benzoylhydrazone 4-oxime	207-807-9	495-73-8	Acute Tox. 3 * Acute Tox. 4 *	H301 H312	GHS06 Dgr	H301 H312			
650-007-00-3	chlordimeform (ISO); N_2-(4-chloro-*o*-tolyl)-N_1,N_1-dimethylformamidine	228-200-5	6164-98-3	Carc. 2 Acute Tox. 4 * Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H351 H312 H302 H400 H410	GHS08 GHS07 GHS09 Wng	H351 H312 H302 H410			
650-008-00-9	drazoxolon (ISO); 4-(2-chlorophenylhydrazone)-3-methyl-5-isoxazolone	227-197-8	5707-69-7	Acute Tox. 3 * Aquatic Acute 1 Aquatic Chronic 1	H301 H400 H410	GHS06 GHS09 Dgr	H301 H410			
650-009-00-4	chlordimeform hydrochloride; *N*'-(4-chloro-*o*-tolyl)-*N*,*N*-dimethylformamidine monohydrochloride; N^2-(4-chloro-*o*-tolyl)-N^1,N^1-dimethylformamidine hydorchloride	243-269-1	19750-95-9	Carc. 2 Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H351 H302 H400 H410	GHS08 GHS07 GHS09 Wng	H351 H302 H410			
650-010-00-X	benzyl violet 4B; α-[4-(4-dimethylamino-α-{4-[ethyl(3-sodiosulphonatobenzyl)amino] phenyl}benzylidene)cyclohexa-2,5-dienylidene(ethyl) ammonio]toluene-3-sulphonate	216-901-9	1694-09-3	Carc. 2	H351	GHS08 Wng	H351			
650-012-00-0	erionite	—	12510-42-8	Carc. 1A	H350	GHS08 Dgr	H350			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
650-013-00-6	asbestos	— — — — — — — —	12001-28-4 132207-32-0 12172-73-5 77536-66-4 77536-68-6 77536-67-5 12001-29-5	Carc. 1A STOT RE 1	H350 H372 **	GHS08 Dgr	H350 H372 **			
650-014-00-1	diethyl 2,4-dihydroxycyclodisiloxane-2,4-diylbis(trimethylene)diphosphonate, tetrasodium salt, reaction products with disodium metasilicate	401-770-4	—	Skin Corr. 1B Acute Tox. 4 *	H314 H302	GHS05 GHS07 Dgr	H314 H302			
650-015-00-7	rosin; colophony	232-475-7 232-484-6 277-299-1	8050-09-7 8052-10-6 73138-82-6	Skin Sens. 1	H317	GHS07 Wng	H317			
650-016-00-2	Mineral wool, with the exception of those specified elsewhere in this Annex; [Man-made vitreous (silicate) fibres with random orientation with alkaline oxide and alkali earth oxide ($Na_2O+K_2O+CaO+MgO+BaO$) content greater than 18 % by weight]	—	—	Carc. 2 Skin Irrit. 2	H351 H315	GHS08 GHS07 Wng	H351 H315			AQR
650-017-00-8	Refractory Ceramic Fibres; Special Purpose Fibres, with the exception of those specified elsewhere in this Annex; [Man-made vitreous (silicate) fibres with random orientation with alkaline oxide and alkali earth oxide ($Na_2O+K_2O+CaO+MgO+BaO$) content less or equal to 18 % by weight]	—	—	Carc. 1B Skin Irrit. 2	H350i H315	GHS08 GHS07 Dgr	H350i H315			A R

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
650-018-00-3	Reaction product of: acetophenone, formaldehyde, cyclohexylamine, methanol and acetic acid	406-230-1	—	Flam. Liq. 3 Carc. 2 Skin Corr. 1B Acute Tox. 4 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H226 H351 H314 H332 H317 H400 H410	GHS02 GHS08 GHS05 GHS07 GHS09 Dgr	H226 H351 H314 H332 H317 H410			
650-031-00-4	bis(4-hydroxy-*N*-methylanilinium) sulphate	200-237-1	55-55-0	Acute Tox. 4 * STOT RE 2 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H302 H373 ** H317 H400 H410	GHS08 GHS07 GHS09 Wng	H302 H373 ** H317 H410			
650-032-00-X	cyproconazole (ISO); (2*RS*,3*RS*;2*RS*,3*SR*)-2-(4-chlorophenyl)-3-cyclopropyl-1-(1*H*—1,2,4-triazol-1-yl) butan-2-ol	—	94361-06-5	Repr. 2 Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H361d *** H302 H400 H410	GHS08 GHS07 GHS09 Wng	H361d *** H302 H410			
650-033-00-5	(*S*)-α-cyano-3-phenoxybenzyl-(*S*)-2-(4-chlorophenyl)-3-methylbutyrate; esfenvalerate	—	66230-04-4	Acute Tox. 3 * Acute Tox. 3 * Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H331 H301 H317 H400 H410	GHS06 GHS09 Dgr	H331 H301 H317 H410			
650-041-00-9	triasulfuron (ISO); 1-[2-(2-chloroethoxy)phenylsulfonyl]-3-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)urea	—	82097-50-5	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			
650-042-00-4	reaction product of: polyethylene-polyamine-(C_{16}-C_{18})-alkylamides with monothio-(C_2)-alkyl phosphonates	417-450-2	—	Eye Irrit. 2 Skin Irrit. 2 Skin Sens. 1 Aquatic Chronic 3	H319 H315 H317 H412	GHS07 Wng	H319 H315 H317 H412			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
650-043-00-X	reaction product of: 3,5-bis-*tert*-butylsalicylic acid and aluminiumsulfate	420-310-3	—	Acute Tox. 4 * Aquatic Acute 1 Aquatic Chronic 1	H302 H400 H410	GHS07 GHS09 Wng	H302 H410			
650-044-00-5	mixed linear and branched C_{14-15} alcohols ethoxylated, reaction product with epichlorohydrin	420-480-9	158570-99-1	Skin Irrit. 2 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H315 H317 H400 H410	GHS07 GHS09 Wng	H315 H317 H410			
650-045-00-0	reaction product of: 1,2,3-propanetricarboxylic acid, 2-hydroxy, diethyl ester, 1-propanol and zirconium tetra-n-propanolate	417-110-3	—	Flam. Liq. 2 Skin Irrit. 2 Eye Dam. 1 Aquatic Chronic 2	H225 H315 H318 H411	GHS02 GHS05 GHS09 Dgr	H225 H315 H318 H411			
650-046-00-6	di(tetramethylammonium)(29*H*,31*H*-phthalocyanin-*N*29,*N*30,*N*31,*N*32)disulfonamide disulfonate, cuprate(2-)complex, derivates	416-180-2	12222-04-7	Acute Tox. 4 * STOT RE 2 * Aquatic Chronic 2	H302 H373 ** H411	GHS08 GHS07 GHS09 Wng	H302 H373 ** H411			
650-047-00-1	dibenzylphenylsulfonium hexafluoroantimonate	417-760-8	134164-24-2	STOT RE 1 Acute Tox. 4 * Eye Dam. 1 Skin Sens. 1 Aquatic Chronic 2	H372 ** H302 H318 H317 H411	GHS08 GHS05 GHS07 GHS09 Dgr	H372 ** H302 H318 H317 H411			
650-048-00-7	reaction product of: borax, hydrogen peroxide, acetic acid anhydride and acetic acid	420-070-1	—	Org. Perox. D **** Acute Tox. 4 * Acute Tox. 4 * Acute Tox. 4 * Skin Corr. 1A Aquatic Acute 1	H242 H332 H312 H302 H314 H400	GHS02 GHS05 GHS07 GHS09 Dgr	H242 H332 H312 H302 H314 H400			

Index No	International Chemical Identification	EC No	CAS No	Classification		Labelling			Specific Conc. Limits, M-factors	Notes
				Hazard Class and Category Code(s)	Hazard statement Code(s)	Pictogram, Signal Word Code(s)	Hazard statement Code(s)	Suppl. Hazard statement Code(s)		
650-049-00-2	2-alkoyloxyethyl hydrogen maleate, where alkoyl represents (by weight) 70 to 85 % unsaturated octadecoyl, 0.5 to 10 % saturated octadecoyl, and 2 to 18 % saturated hexadecoyl	417-960-5	—	Skin Irrit. 2 Eye Dam. 1 Skin Sens. 1 Aquatic Acute 1 Aquatic Chronic 1	H315 H318 H317 H400 H410	GHS05 GHS07 GHS09 Dgr	H315 H318 H317 H410			
650-050-00-8	reaction mass of: 1-methyl-3-hydroxypropyl 3,5-[1,1-dimethylethyl]-4-hydroxydihydrocinnamate and/or 3-hydroxybutyl 3,5-[1,1-dimethylethyl]-4-hydroxydihydrocinnamate; 1,3-butanediol bis[3-(3'-(1,1-dimethylethyl) 4'-hydroxy-phenyl)propionate] isomers; 1,3-butanediol bis[3-(3',5'-(1,1-dimethylethyl)-4'-hydroxyphenyl)propionate] isomers	423-600-8	—	Aquatic Chronic 2	H411	GHS09	H411			
650-055-00-5	silver sodium zirconium hydrogenphosphate	422-570-3	155925-27-2	Aquatic Acute 1 Aquatic Chronic 1	H400 H410	GHS09 Wng	H410			

Table 3.2

The list of harmonised classification and labelling of hazardous substances from Annex I to Directive 67/548/EEC

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
001-001-00-9	hydrogen	215-605-7	1333-74-0	F+; R12	F+ R: 12 S: (2-)9-16-33		
001-002-00-4	aluminium lithium hydride	240-877-9	16853-85-3	F; R15	F R: 15 S: (2-)7/8-24/25-43		
001-003-00-X	sodium hydride	231-587-3	7646-69-7	F; R15	F R: 15 S: (2-)7/8-24/25-43		
001-004-00-5	calcium hydride	232-189-2	7789-78-8	F; R15	F R: 15 S: (2-)7/8-24/25-43		
003-001-00-4	lithium	231-102-5	7439-93-2	F; R15 R14 C; R34	F; C R: 14/15-34 S: (1/2-)8-43-45		
003-002-00-X	n-hexyllithium	404-950-0	21369-64-2	F; R15-17 R14 C; R35	F; C R: 14/15-17-35 S: (1/2-)6-16-26-30-36/37/39-43-45		
004-001-00-7	beryllium	231-150-7	7440-41-7	Carc. Cat. 2; R49 T+; R26 T; R25-48/23 Xi; R36/37/38 R43	T+ R: 49-25-26-36/37/38-43-48/23 S: 53-45		E

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
004-002-00-2	beryllium compounds with the exception of aluminium beryllium silicates, and with those specified elsewhere in this Annex	—	—	Carc. Cat. 2; R49 T+; R26 T; R25-48/23 Xi; R36/37/38 R43 N; R51-53	T+; N R: 49-25-26-36/37/38-43-48/23-51/53 S: 53-45-61		AE
004-003-00-8	beryllium oxide	215-133-1	1304-56-9	Carc. Cat. 2; R49 T+; R26 T; R25-48/23 Xi; R36/37/38 R43	T+ R: 49-25-26-36/37/38-43-48/23 S: 53-45		E
005-001-00-X	boron trifluoride	231-569-5	7637-07-2	R14 T+; R26 C; R35	T+; C R: 14-26-35 S: (1/2-)9-26-28-36/37/39-45		
005-002-00-5	boron trichloride	233-658-4	10294-34-5	R14 T+; R26/28 C; R34	T+ R: 14-26/28-34 S: (1/2-)9-26-28-36/37/39-45		
005-003-00-0	boron tribromide	233-657-9	10294-33-4	R14 T+; R26/28 C; R35	T+; C R: 14-26/28-35 S: (1/2-)9-26-28-36/37/39-45		
005-004-00-6	trialkylboranes	—	—	F; R17 C; R34	F; C R: 17-34 S: (1/2-)7-23-26-36/37/39-43-45		A
005-005-00-1	trimethyl borate	204-468-9	121-43-7	R10 Xn; R21	Xn R: 10-21 S: (2-)23-25		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
005-006-00-7	dibutyltin hydrogen borate	401-040-5	75113-37-0	T; R48/25 Xn; R21/22 Xi; R41 R43 N; R50-53	T; N R: 21/22-41-43-48/25-50/53 S: (1/2-)22-26-36/37-45-60-61		
005-009-00-3	tetrabutylammonium butyltriphenylborate	418-080-4	120307-06-4	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-56-61		
005-010-00-9	*N*,*N*-dimethylanilinium tetrakis(pentafluorophenyl)borate	422-050-6	118612-00-3	Carc. Cat. 3; R40 Xn; R22 Xi; R38-41	Xn R: 22-38-40-41 S: (2-)22-26-36/37/39		
005-012-00-X	diethyl{4-[1,5,5-tris(4-diethylaminophenyl)penta-2,4-dienylidene]cyclohexa-2,5-dienylidene}ammonium butyltriphenylborate	418-070-1	141714-54-7	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
006-001-00-2	carbon monoxide	211-128-3	630-08-0	F+; R12 Repr. Cat. 1; R61 T; R23-48/23	F+; T R: 61-12-23-48/23 S: 53-45		E
006-002-00-8	phosgene; carbonyl chloride	200-870-3	75-44-5	T+; R26 C; R34	T+ R: 26-34 S: (1/2-)9-26-36/37/39-45		
006-003-00-3	carbon disulphide	200-843-6	75-15-0	F; R11 Repr. Cat. 3; R62-63 T; R48/23 Xi; R36/38	F; T R: 11-36/38-48/23-62-63 S: (1/2-)16-33-36/37-45	Repr. Cat. 3; R62-63: C ≥ 1 % T; R48/23: C ≥ 1 % Xn; R48/20: 0,2 % ≤ C < 1 %	
006-004-00-9	calcium carbide	200-848-3	75-20-7	F; R15	F R: 15 S: (2-)8-43		
006-005-00-4	thiram (ISO); tetramethylthiuram disulphide	205-286-2	137-26-8	Xn; R20/22-48/22 Xi; R36/38 R43 N; R50-53	Xn; N R: 20/22-36/38-43-48/22-50/53 S: (2-)26-36/37-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
006-006-00-X	hydrogen cyanide; hydrocyanic acid	200-821-6	74-90-8	F+; R12 T+; R26 N; R50-53	F+; T+; N R: 12-26-50/53 S: (1/2-)7/9-16-36/37-38-45-60-61		
006-006-01-7	hydrogen cyanide ...%; hydrocyanic acid ...%	200-821-6	74-90-8	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)7/9-36/37-38-45-60-61		B
006-007-00-5	salts of hydrogen cyanide with the exception of complex cyanides such as ferrocyanides, ferricyanides and mercuric oxycyanide	—	—	T+; R26/27/28 R32 N; R50-53	T+; N R: 26/27/28-32-50/53 S: (1/2-)7-28-29-45-60-61		A
006-008-00-0	antu (ISO); 1-(1-naphthyl)-2-thiourea	201-706-3	86-88-4	T+; R28 Carc. Cat. 3; R40	T+ R: 28-40 S: (1/2-)25-36/37-45		
006-009-00-6	1-isopropyl-3-methylpyrazol-5-yl dimethylcarbamate; isolan	204-318-2	119-38-0	T+; R27/28	T+ R: 27/28 S: (1/2-)28-36/37/39-45		
006-010-00-1	5,5-dimethyl-3-oxocyclohex-1-enyl dimethylcarbamate 5,5-dimethyldihydroresorcinol dimethylcarbamate; dimetan	204-525-8	122-15-6	T; R25	T R: 25 S: (1/2-)36/37-45		
006-011-00-7	carbaryl (ISO); 1-naphthyl methylcarbamate	200-555-0	63-25-2	Carc. Cat. 3; R40 Xn; R22 N; R50	Xn; N R: 22-40-50 S: (2-)22-24-36/37-46-61		
006-012-00-2	ziram (ISO); zinc bis dimethyldithiocarbamate	205-288-3	137-30-4	T+; R26 Xn; R22-48/22 Xi; R37-41 R43 N; R50-53	T+; N R: 22-26-37-41-43-48/22-50/53 S: (1/2-)22-26-28-36/37/39-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
006-013-00-8	metam-sodium (ISO); sodium methyldithiocarbamate	205-293-0	137-42-8	Xn; R22 R31 C; R34 R43 N; R50-53	C; N R: 22-31-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
006-014-00-3	nabam (ISO); disodium ethylenebis(*N,N'*-dithiocarbamate)	205-547-0	142-59-6	Xn; R22 Xi; R37 R43 N; R50-53	Xn; N R: 22-37-43-50/53 S: (2-)8-24/25-46-60-61		
006-015-00-9	diuron (ISO); 3-(3,4-dichlorophenyl)-1,1-dimethylurea	206-354-4	330-54-1	Carc. Cat. 3; R40 Xn; R22-48/22 N; R50-53	Xn; N R: 22-40-48/22-50/53 S: (2-)13-22-23-37-46-60-61		
006-016-00-4	propoxur (ISO); 2-isopropyloxyphenyl *N*-methylcarbamate; 2-isopropoxyphenyl methylcarbamate	204-043-8	114-26-1	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)37-45-60-61		
006-017-00-X	aldicarb (ISO); 2-methyl-2-(methylthio)propanal-*O*-(*N*-methylcarbamoyl) oxime	204-123-2	116-06-3	T+; R26/28 T; R24 N; R50-53	T+; N R: 24-26/28-50/53 S: (1/2-)22-36/37-45-60-61		
006-018-00-5	aminocarb (ISO); 4-dimethylamino-3-tolyl methylcarbamate	217-990-7	2032-59-9	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61		
006-019-00-0	di-allate (ISO); *S*-(2,3-dichloroallyl)-*N,N*-diisopropylthiocarbamate	218-961-1	2303-16-4	Carc. Cat. 3; R40 Xn; R22 N; R50-53	Xn; N R: 22-40-50/53 S: (2-)25-36/37-60-61		
006-020-00-6	barban (ISO); 4-chlorbut-2-ynyl *N*-(3-chlorophenyl)carbamate	202-930-4	101-27-9	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-36/37-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
006-021-00-1	linuron (ISO); 3-(3,4-dichlorophenyl)-1-methoxy-1-methylurea	206-356-5	330-55-2	Repr. Cat. 2; R61 Repr. Cat. 3; R62 Carc. Cat. 3; R40 Xn; R22-48/22 N; R50-53	T; N R: 61-22-40-48/22-62-50/53 S: 53-45-60-61		E
006-022-00-7	decarbofuran (ISO); 2,3-dihydro-2-methylbenzofuran-7-yl methylcarbamate	—	1563-67-3	T; R23/24/25	T R: 23/24/25 S: (1/2-)13-36/37-45		
006-023-00-2	mercaptodimethur (ISO); methiocarb (ISO); 3,5-dimethyl-4-methylthiophenyl *N*-methylcarbamate	217-991-2	2032-65-7	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)22-37-45-60-61		
006-024-00-8	proxan-sodium (ISO); sodium *O*-isopropyldithiocarbonate	205-443-5	140-93-2	Xn; R22 Xi; R38 N; R51-53	Xn; N R: 22-38-51/53 S: (2-)13-61		
006-025-00-3	allethrin; (*RS*)-3-allyl-2-methyl-4-oxocyclopent-2-enyl (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-(2-methylprop-1-enyl) cyclopropanecarboxylate; bioallethrin; (*RS*)-3-allyl-2-methyl-4-oxocyclopent-2-enyl (1*R*,3*R*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate; [1] S-bioallethrin; (*S*)-3-allyl-2-methyl-4-oxocyclopent-2-enyl (1*R*,3*R*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate; [2] esbiothrin; (*RS*)-3-allyl-2-methyl-4-oxocyclopent-2-enyl (1*R*,3*R*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate [3]	209-542-4 [1] 249-013-5 [2] - [3]	584-79-2 [1] 28434-00-6 [2] 84030-86-4 [3]	Xn; R20/22 N; R50-53	Xn; N R: 20/22-50/53 S: (2-)36-60-61		C
006-026-00-9	carbofuran (ISO); 2,3-dihydro-2,2-dimethylbenzofuran-7-yl *N*-methylcarbamate	216-353-0	1563-66-2	T+; R26/28 N; R50-53	T+; N R: 26/28-50/53 S: (1/2-)36/37-45-60-61		

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006-028-00-X	dinobuton (ISO); 2-(1-methylpropyl)-4,6-dinitrophenyl isopropyl carbonate	213-546-1	973-21-7	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)37-45-60-61		
006-029-00-5	dioxacarb (ISO); 2-(1,3-dioxolan-2-yl)phenyl *N*-methylcarbamate	230-253-4	6988-21-2	T; R25 N; R51-53	T; N R: 25-51/53 S: (1/2-)37-45-61		
006-030-00-0	EPTC (ISO); *S*-ethyl dipropylthiocarbamate	212-073-8	759-94-4	Xn; R22	Xn R: 22 S: (2-)23		
006-031-00-6	formetanate (ISO); 3-[(EZ)-dimethylaminomethyleneamino]phenyl methylcarbamate	244-879-0	22259-30-9	T+; R26/28 R43 N; R50-53	T+; N R: 26/28-43-50/53 S: (1/2-)24-28-37/39-45-60-61		
006-032-00-1	monolinuron (ISO); 3-(4-chlorophenyl)-1-methoxy-1-methylurea	217-129-5	1746-81-2	Xn; R22-48/22 N; R50-53	Xn; N R: 22-48/22-50/53 S: (2-)22-60-61		
006-033-00-7	metoxuron (ISO); 3-(3-chloro-4-methoxyphenyl)-1,1-dimethylurea	243-433-2	19937-59-8	N; R50-53	N R: 50/53 S: 60-61		
006-034-00-2	pebulate (ISO); *N*-butyl-*N*-ethyl-*S*-propylthiocarbamate	214-215-4	1114-71-2	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)23-61		
006-035-00-8	pirimicarb (ISO); 5,6-dimethyl-2-dimethylamino-pyrimidin-4-yl *N*,*N*-dimethylcarbamate	245-430-1	23103-98-2	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)22-37-45-60-61		
006-036-00-3	benzthiazuron (ISO); 1-benzothiazol-2-yl-3-methylurea	217-685-9	1929-88-0	Xn; R22	Xn R: 22 S: (2-)24/25		
006-037-00-9	promecarb (ISO); 3-isopropyl-5-methylphenyl *N*-methylcarbamate	220-113-0	2631-37-0	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)24-37-45-60-61		

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006-038-00-4	sulfallate (ISO); 2-chloroallyl *N*,*N*-dimethyldithiocarbamate	202-388-9	95-06-7	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		E
006-039-00-X	tri-allate (ISO); *S*-2,3,3-trichloroallyl diisopropylthiocarbamate	218-962-7	2303-17-5	Xn; R22-48/22 R43 N; R50-53	Xn; N R: 22-43-48/22-50/53 S: (2-)24-37-60-61		
006-040-00-5	3-methylpyrazol-5-yl-dimethylcarbamate; monometilan	—	2532-43-6	T; R23/24/25	T R: 23/24/25 S: (1/2-)13-45		
006-041-00-0	dimethylcarbamoyl chloride	201-208-6	79-44-7	Carc. Cat. 2; R45 T; R23 Xn; R22 Xi; R36/37/38	T R: 45-22-23-36/37/38 S: 53-45	Carc. Cat. 2; R45: C ≥ 0,001 %	E
006-042-00-6	monuron (ISO); 3-(4-chlorophenyl)-1,1-dimethylurea	205-766-1	150-68-5	Carc. Cat. 3; R40 Xn; R22 N; R50-53	Xn; N R: 22-40-50/53 S: (2-)36/37-60-61		
006-043-00-1	3-(4-chlorophenyl)-1,1-dimethyluronium trichloroacetate; monuron-TCA	—	140-41-0	Xi; R36/38 Carc. Cat. 3; R40 N; R50-53	Xn; N R: 36/38-40-50/53 S: (2-)36/37-60-61		
006-044-00-7	isoproturon (ISO); 3-(4-isopropylphenyl)-1,1-dimethylurea	251-835-4	34123-59-6	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
006-045-00-2	methomyl (ISO); 1-(methylthio)ethylideneamino *N*-methylcarbamate	240-815-0	16752-77-5	T+;R28 N; R50-53	T+; N R: 28-50/53 S: (1/2-)22-36/37-45-60-61		
006-046-00-8	bendiocarb (ISO); 2,2-dimethyl-1,3-benzodioxol-4-yl N-methylcarbamate	245-216-8	22781-23-3	T; R23/25 Xn; R21 N; R50-53	T; N R: 21-23/25-50/53 S: (1/2-)22-36/37-45-60-61		

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006-047-00-3	bufencarb (ISO); reaction mass of 3-(1-methylbutyl)phenyl *N*-methylcarbamate and 3-(1-ethylpropyl)phenyl *N*-methylcarbamate	—	8065-36-9	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61		
006-048-00-9	ethiofencarb (ISO); 2-(ethylthiomethyl)phenyl *N*-methylcarbamate	249-981-9	29973-13-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
006-049-00-4	dixanthogen; *O*,*O*-diethyl dithiobis(thioformate)	207-944-4	502-55-6	Xn; R22	Xn R: 22 S: (2-)24		
006-050-00-X	1,1-dimethyl-3-phenyluronium trichloroacetate; fenuron-TCA	—	4482-55-7	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)60-61		
006-051-00-5	ferbam (ISO); iron tris(dimethyldithiocarbamate)	238-484-2	14484-64-1	Xi; R36/37/38 N; R50-53	Xi; N R: 36/37/38-50/53 S: (2-)60-61		
006-052-00-0	formetanate hydrochloride; 3-(*N*,*N*-dimethylaminomethyleneamino)phenyl *N*-methylcarbamate	245-656-0	23422-53-9	T+; R26/28 R43 N; R50-53	T+; N R: 26/28-43-50/53 S: (1/2-)24-28-37/39-45-60-61		
006-053-00-6	isoprocarb (ISO); 2-isopropylphenyl *N*-methylcarbamate	220-114-6	2631-40-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
006-054-00-1	mexacarbate (ISO); 3,5-dimethyl-4-dimethylaminophenyl *N*-methylcarbamate	206-249-3	315-18-4	T+; R28 Xn; R21 N; R50-53	T+; N R: 21-28-50/53 S: (1/2-)36/37-45-60-61		
006-055-00-7	xylylcarb (ISO); 3,4-dimethylphenyl N-methylcarbamate; 3,4-xylyl methylcarbamate; MPMC	219-364-9	2425-10-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
006-056-00-2	metolcarb (ISO); *m*-tolyl methylcarbamate; MTMC	214-446-0	1129-41-5	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		

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006-057-00-8	nitrapyrin (ISO); 2-chloro-6-trichloromethylpyridine	217-682-2	1929-82-4	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)24-61		
006-058-00-3	noruron (ISO); 1,1-dimethyl-3-(perhydro-4,7-methanoinden-5-yl)urea	—	2163-79-3	Xn; R22	Xn R: 22 S: (2-)		
006-059-00-9	oxamyl (ISO); *N*',*N*'-dimethylcarbamoyl(methylthio)methylenamine *N*-methylcarbamate;	245-445-3	23135-22-0	T+; R26/28 Xn; R21 N; R51-53	T+; N R: 21-26/28-51/53 S: (1/2-)36/37-45-61		
006-060-00-4	oxycarboxin (ISO); 2,3-dihydro-6-methyl-5-(*N*-phenylcarbamoyl)-1,4-oxothiine 4,4-dioxide	226-066-2	5259-88-1	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
006-061-00-X	*S*-ethyl *N*-(dimethylaminopropyl)thiocarbamatehydrochloride; prothiocarb hydrochloride	243-193-9	19622-19-6	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
006-062-00-5	methyl 3,4-dichlorophenylcarbanilate; SWEP.	—	1918-18-9	Xn; R22	Xn R: 22 S: (2-)		
006-063-00-0	thiobencarb (ISO); *S*-4-chlorobenzyl diethylthiocarbamate;	248-924-5	28249-77-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
006-064-00-6	thiofanox (ISO); 3,3-dimethyl-1-(methylthio)butanone-*O*-(*N*-methylcarbamoyl)oxime;	254-346-4	39196-18-4	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)27-36/37-45-60-61		
006-065-00-1	3-chloro-6-cyano-bicyclo(2,2,1)heptan-2-one-*O*-(*N*-methylcarbamoyl)oxime; triamid	—	15271-41-7	T+; R28 T; R24 N; R51-53	T+; N R: 24-28-51/53 S: (1/2-)28-36/37-45-61		
006-066-00-7	vernolate (ISO); *S*-propyl dipropylthiocarbamate;	217-681-7	1929-77-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		

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006-067-00-2	XMC; 3,5-xylyl methylcarbamate	—	2655-14-3	Xn; R22	Xn R: 22 S: (2-)		
006-068-00-8	diazomethane	206-382-7	334-88-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		
006-069-00-3	thiophanate-methyl (ISO); 1,2-di-(3-methoxycarbonyl-2-thioureido)benzene	245-740-7	23564-05-8	Muta. Cat. 3; R68 Xn; R20 R43 N; R50-53	Xn; N R: 20-43-50/53-68 S: (2-)36/37-46-60-61		
006-070-00-9	furmecyclox (ISO); *N*-cyclohexyl-*N*-methoxy-2,5-dimethyl-3-furamide;	262-302-0	60568-05-0	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		
006-071-00-4	cyclooct-4-en-1-yl methyl carbonate	401-620-8	87731-18-8	R43	Xi R: 43 S: (2-)24-37		
006-072-00-X	prosulfocarb(ISO); *S*-benzyl *N*,*N*-dipropylthiocarbamate;	401-730-6	52888-80-9	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)24-37-61		
006-073-00-5	3-(dimethylamino)propylurea	401-950-2	31506-43-1	Xi; R41	Xi R: 41 S: (2-)26-39		
006-074-00-0	2-(3-(prop-1-en-2-yl)phenyl)prop-2-yl isocyanate	402-440-2	2094-99-7	T+; R26 C; R34 Xn; R48/20 R42/43 N; R50-53	T+; N R: 26-34-42/43-48/20-50/53 S: (1/2-)7-15-28-36/37/39-38-45-60-61		
006-076-00-1	mancozeb (ISO)	—	8018-01-7	Xi; R37 R43	Xi R: 37-43 S: (2-)8-24/25-46		
006-077-00-7	maneb (ISO)	235-654-8	12427-38-2	Xi; R37 R43	Xi R: 37-43 S: (2-)8-24/25-46		

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006-078-00-2	zineb (ISO); zinc ethylenebis(dithiocarbamate) (polymeric)	235-180-1	12122-67-7	Xi; R37 R43	Xi R: 37-43 S: (2-)8-24/25-46		
006-079-00-8	disulfiram; tetraethylthiuramdisulfide	202-607-8	97-77-8	Xn; R22-48/22 R43 N; R50-53	Xn; N R: 22-43-48/22-50/53 S: (2-)24-37-60-61		
006-080-00-3	tetramethylthiuram monosulphide	202-605-7	97-74-5	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)24-26-37-61		
006-081-00-9	zinc bis(dibutyldithiocarbamate)	205-232-8	136-23-2	Xi; R36/37/38 R43 N; R50-53	Xi; N R: 36/37/38-43-50/53 S: (2-)24-37-60-61		
006-082-00-4	zinc bis(diethyldithiocarbamate)	238-270-9	14324-55-1	Xn; R22 Xi; R36/37/38 R43 N; R50-53	Xn; N R: 22-36/37/38-43-50/53 S: (2-)24-37-60-61		
006-083-00-X	butocarboxim (ISO); 3-(methylthio)-2-butanone *O*-[(methylamino)carbonyl]oxime	252-139-3	34681-10-2	R10 T; R23/24/25 Xi; R36 N; R50-53	T; N R: 10-23/24/25-36-50/53 S: (1/2-)36/37-45-60-61		
006-084-00-5	carbosulfan (ISO); 2,3-dihydro-2,2-dimethyl-7-benzofuryl [(dibutylamino)thio] methylcarbamate;	259-565-9	55285-14-8	T; R23/25 R43 N; R50-53	T; N R: 23/25-43-50/53 S: (1/2-)24-37-38-45-60-61		
006-085-00-0	fenobucarb (ISO); 2-butylphenyl methylcarbamate;	223-188-8	3766-81-2	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
006-086-00-6	ethyl [2-(4-phenoxyphenoxy)ethyl]carbamate; fenoxycarb	276-696-7	72490-01-8	N; R50-53	N R: 50/53 S: 60-61		

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006-087-00-1	2,3-dihydro-2,2-dimethyl-7-benzofuryl 2,4-dimethyl-6-oxa-5-oxo-3-thia-2,4-diazadecanoate; furathiocarb	265-974-3	65907-30-4	T+; R26 T; R25 Xn; R48/22 Xi; R36/38 R43 N; R50-53	T+; N R: 25-26-36/38-43-48/22-50/53 S: (1/2-)28-36/37-38-45-60-61		
006-088-00-7	benfuracarb; ethyl *N*-[2,3-dihydro-2,2-dimethylbenzofuran-7-yloxycarbonyl(methyl)aminothio]-*N*-isopropyl- β-alaninate	—	82560-54-1	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)36/37-45-60-61		
006-089-00-2	chlorine dioxide	233-162-8	10049-04-4	O; R8 R6 T+; R26 C; R34 N; R50	O; T+; N R: 6-8-26-34-50 S: (1/2-)23-26-28-36/37/39-38-45-61	N; R50: C ≥ 0,02 %	5
006-089-01-X	chlorine dioxide … %	233-162-8	10049-04-4	T; R25 C; R34 N; R50	T; N R: 25-34-50 S: (1/2-)23-26-28-36/37/39-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 3 % ≤ C < 10 % Xi; R36: 0,3 % ≤ C < 3 % N; R50: C ≥ 3 %	B
006-090-00-8	2-(3-iodoprop-2-yn-1-yloxy)ethyl phenylcarbamate	408-010-0	88558-41-2	Xn; R20 Xi; R41 R52-53	Xn R: 20-41-52/53 S: (2-)22-26-39-61		
007-001-00-5	ammonia, anhydrous	231-635-3	7664-41-7	R10 ⊗ T; R23 C; R34 N; R50	T; N R: 10-23-34-50 S: (1/2-)9-16-26-36/37/39-45-61		
007-001-01-2	ammonia ....%	215-647-6	1336-21-6	C; R34 N; R50	C; N R: 34-50 S: (1/2-)26-36/37/39-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	B
007-002-00-0	nitrogen dioxide; [1] dinitrogen tetraoxide [2]	233-272-6 [1] 234-126-4 [2]	10102-44-0 [1] 10544-72-6 [2]	T+; R26 C; R34	T+ R: 26-34 S: (1/2-)9-26-28-36/37/39-45	T+; R26: C ≥ 10 % T; R23: 1 % ≤ C < 10 % Xn; R20: 0,1 % ≤ C < 1 %	5

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007-003-00-6	chlormequat chloride (ISO); 2-chloroethyltrimethylammonium chloride	213-666-4	999-81-5	Xn; R21/22	Xn R: 21/22 S: (2-)36/37		
007-004-00-1	nitric acid ... %	231-714-2	7697-37-2	O; R8 C; R35	O; C R: 8-35 S: (1/2-)23-26-36-45	C; R35: C ≥ 20 % C; R34: 5 % ≤ C < 20 % Footnote: O; R8: C ≥ 70 %	B
007-006-00-2	ethyl nitrite	203-722-6	109-95-5	E; R2 Xn; R20/21/22	E; Xn R: 2-20/21/22 S: (2-)		
007-007-00-8	ethyl nitrate	210-903-3	625-58-1	E; R2 ⊗	E R: 2 S: (2-)23-24/25		
007-008-00-3	hydrazine	206-114-9	302-01-2	R10 Carc. Cat. 2; R45 T; R23/24/25 C; R34 R43 N; R50-53	T; N R: 45-10-23/24/25-34-43-50/53 S: 53-45-60-61	C; R34: C ≥ 10 % Xi; R36/38: 3 % ≤ C < 10 %	E
007-009-00-9	dicyclohexylammonium nitrite	221-515-9	3129-91-7	Xn; R20/22	Xn R: 20/22 S: (2-)15-41	Xn; R20/22: C ≥ 10 %	
007-010-00-4	sodium nitrite	231-555-9	7632-00-0	O; R8 T; R25 N; R50	O; T; N R: 8-25-50 S: (1/2-)45-61	T; R25: C ≥ 5 % Xn; R22: 1 % ≤ C < 5 %	
007-011-00-X	potassium nitrite	231-832-4	7758-09-0	O; R8 T; R25 N; R50	O; T; N R: 8-25-50 S: (1/2-)45-61	T; R25: C ≥ 5 % Xn; R22: 1 % ≤ C < 5 %	
007-012-00-5	*N,N*-dimethylhydrazine	200-316-0	57-14-7	F; R11 Carc. Cat. 2; R45 T; R23/25 C; R34 N; R51-53	F; T; N R: 45-11-23/25-34-51/53 S: 53-45-61		E

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007-013-00-0	1,2-dimethylhydrazine	—	540-73-8	Carc. Cat. 2; R45 T; R23/24/25 N; R51-53	T; N R: 45-23/24/25-51/53 S: 53-45-61	Carc. Cat. 2; R45: C ≥ 0,01 %	E
007-014-00-6	salts of hydrazine	—	—	Carc. Cat. 2; R45 T; R23/24/25 R43 N; R50-53	T; N R: 45-23/24/25-43-50/53 S: 53-45-60-61		AE
007-015-00-1	*O*-ethylhydroxylamine	402-030-3	624-86-2	F; R11 T; R23/24/25-48/23 Xi; R36 R43 N; R50	F; T; N R: 11-23/24/25-36-43-48/23-50 S: (1/2-)16-26-36/37/39-45-60-61		
007-016-00-7	butyl nitrite	208-862-1	544-16-1	F; R11 T; R23/25	F; T R: 11-23/25 S: (1/2-)16-24-45		
007-017-00-2	isobutyl nitrite	208-819-7	542-56-3	F; R11 Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R20/22	F; T R: 11-20/22-45-68 S: 53-45		E
007-018-00-8	*sec*-butyl nitrite	213-104-8	924-43-6	F; R11 Xn; R20/22	F; Xn R: 11-20/22 S: (2-)16-24-46		
007-019-00-3	*tert*-butyl nitrite	208-757-0	540-80-7	F; R11 Xn; R20/22	F; Xn R: 11-20/22 S: (2-)16-24-46		
007-020-00-9	pentyl nitrite; [1] 'amyl nitrite', mixed isomers [2]	207-332-7 [1] 203-770-8 [2]	463-04-7 [1] 110-46-3 [2]	F; R11 Xn; R20/22	F; Xn R: 11-20/22 S: (2-)16-24-46		

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007-021-00-4	hydrazobenzene; 1,2-diphenylhydrazine	204-563-5	122-66-7	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		E
007-022-00-X	hydrazine bis(3-carboxy-4-hydroxybenzensulfonate)	405-030-1	—	Carc. Cat. 2; R45 Xn; R22 C; R34 R43 R52-53	T R: 45-22-34-43-52/53 S: 53-45-61		E
007-023-00-5	sodium 3,5-bis(3-(2,4-di-tert-pentylphenoxy)propylcarbamoyl)benzenesulfonate	405-510-0	—	Xi; R38 R43	Xi R: 38-43 S: (2-)24-37		
007-024-00-0	2-(decylthio)ethylammonium chloride	405-640-8	36362-09-1	Xn; R48/22 Xi; R38-41 N; R50-53	Xn; N R: 38-41-48/22-50/53 S: (2-)26-36/37/39-60-61		
007-025-00-6	(4-hydrazinophenyl)-*N*-methylmethanesulfonamide hydrochloride	406-090-1	81880-96-8	Muta. Cat. 3; R68 T; R25-48/25 R43 N; R50-53	T; N R: 25-43-48/25-68-50/53 S: (1/2-)22-36/37/39-45-60-61		
007-026-00-1	oxo-((2,2,6,6-tetramethylpiperidin-4-yl)amino)carbonylacetohydrazide	413-230-5	122035-71-6	Xi; R41 R43	Xi R: 41-43 S: (2-)8-22-24-26-30-37/39		
007-027-00-7	1,6-bis(3,3-bis((1-methylpentylidenimino)propyl)ureido)hexane	420-190-2	771478-66-1	C; R34 Xn; R21/22-48/21 R43 N; R50-53	C; N R: 21/22-34-43-48/21-50/53 S: (1/2-)7-26-36/37/39-45-60-61		
008-001-00-8	oxygen	231-956-9	7782-44-7	O; R8	O R: 8 S: (2-)17		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
008-003-00-9	hydrogen peroxide solution ... %	231-765-0	7722-84-1	R5 O; R8 C; R35 Xn; R20/22	O; C R: 5-8-20/22-35 S: (1/2-)17-26-28-36/37/39-45	Xn; R20: C ≥ 50 % Xn; R22: C ≥ 8 % C; R35: C ≥ 70 % C; R34: 50 % ≤ C < 70 % Xi; R37/38: 35 % ≤ C < 50 % Xi; R41: 8 % ≤ C < 50 % Xi; R36: 5 % ≤ C < 8 % Footnote: O; R8: C ≥ 50 % R5: C ≥ 70 %	B
009-001-00-0	fluorine	231-954-8	7782-41-4	R7 ⊗ T+; R26 C; R35	T+; C R: 7-26-35 S: (1/2-)9-26-36/37/39-45		
009-002-00-6	hydrogen fluoride	231-634-8	7664-39-3	T+; R26/27/28 C; R35	T+; C R: 26/27/28-35 S: (1/2-)7/9-26-36/37/39-45		
009-003-00-1	hydrofluoric acid ... %	231-634-8	7664-39-3	T+; R26/27/28 C; R35	T+; C R: 26/27/28-35 S: (1/2-)7/9-26-36/37-45	C; R35: C ≥ 7 % C; R34: 1 % ≤ C < 7 % Xi; R36: 0,1 % ≤ C < 1 %	B
009-004-00-7	sodium fluoride	231-667-8	7681-49-4	T; R25 Xi; R36/38 R32	T R: 25-32-36/38 S: (1/2-)22-36-45		
009-005-00-2	potassium fluoride	232-151-5	7789-23-3	T; R23/24/25	T R: 23/24/25 S: (1/2-)26-45		
009-006-00-8	ammonium fluoride	235-185-9	12125-01-8	T; R23/24/25	T R: 23/24/25 S: (1/2-)26-45		
009-007-00-3	sodium bifluoride; sodium hydrogen difluoride	215-608-3	1333-83-1	T; R25 C; R34	T; C R: 25-34 S: (1/2-)22-26-37-45	T; R25: C ≥ 10 % Xn; R22: 1 % ≤ C < 10 % C; R34: C ≥ 1 % Xi; R36/38: 0,1 % ≤ C < 1 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
009-008-00-9	potassium bifluoride; potassium hydrogen difluoride	232-156-2	7789-29-9	T; R25 C; R34	T; C R: 25-34 S: (1/2-)22-26-37-45	T; R25: C ≥ 10 % Xn; R22: 1 % ≤ C < 10 % C; R34: C ≥ 1 % Xi; R36/38: 0,1 % ≤ C < 1 %	
009-009-00-4	ammonium bifluoride; ammonium hydrogen difluoride	215-676-4	1341-49-7	T; R25 C; R34	T; C R: 25-34 S: (1/2-)22-26-37-45	T; R25: C ≥ 10 % Xn; R22: 1 % ≤ C < 10 % C; R34: C ≥ 1 % Xi; R36/38: 0,1 % ≤ C < 1 %	
009-010-00-X	fluoroboric acid ... %	240-898-3	16872-11-0	C; R34	C R: 34 S: (1/2-)26-27-45	C; R34: C ≥ 25 % Xi; R36/38: 10 % ≤ C < 25 %	B
009-011-00-5	fluorosilicic acid ... %	241-034-8	16961-83-4	C; R34	C R: 34 S: (1/2-)26-27-45		B
009-012-00-0	alkali fluorosilicates(Na); [1] alkali fluorosilicates(K); [2] alkali fluorosilicates(NH4) [3]	240-934-8 [1] 240-896-2 [2] 240-968-3 [3]	16893-85-9 [1] 16871-90-2 [2] 16919-19-0 [3]	T; R23/24/25	T R: 23/24/25 S: (1/2-)26-45	T; R23/24/25: C ≥ 10 % Xn; R20/21/22: 1 % ≤ C < 10 %	A
009-013-00-6	fluorosilicates, with the exception of those specified elsewhere in this annex	—	—	Xn; R22	Xn R: 22 S: (2-)13-24/25	Xn; R22: C ≥ 10 %	A
009-014-00-1	lead hexafluorosilicate	247-278-1	25808-74-6	Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R20/22 R33 N; R50-53	T; N R: 61-62-20/22-33-50/53 S: 53-45-60-61		E 1
009-015-00-7	sulphuryl difluoride	220-281-5	2699-79-8	T; R23 Xn; R48/20 N; R50	T; N R: 23-48/20-50 S: (1/2-)45-63-60-61		

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009-016-00-2	trisodium hexafluoroaluminate; cryolite	237-410-6 239-148-8	13775-53-6 15096-52-3	T; R48/23/25 Xn; R20/22 N; R51-53	T; N R: 20/22-48/23/25-51/53 S: (1/2-)22-37-45-61		C
009-017-00-8	potassium mu-fluoro-bis(triethylaluminium)	400-040-2	12091-08-6	F; R11-14/15 C; R35 Xn; R20	F; C R: 11-14/15-20-35 S: (1/2-)16-30-36/39-43-45		
009-018-00-3	magnesium hexafluorosilicate	241-022-2	16949-65-8	T; R25	T R: 25 S: (1/2-)24/25-45	T; R25: C ≥ 10 % Xn; R22: 1 % ≤ C < 10 %	
011-001-00-0	sodium	231-132-9	7440-23-5	F; 15 R14 C; R34	F; C R: 14/15-34 S: (1/2-)5-8-43-45		
011-002-00-6	sodium hydroxide; caustic soda	215-185-5	1310-73-2	C; R35	C R: 35 S: (1/2-)26-37/39-45	C; R35: C ≥ 5 % C; R34: 2 % ≤ C < 5 % Xi; R36/38: 0,5 % ≤ C < 2 %	
011-003-00-1	sodium peroxide	215-209-4	1313-60-6	O; R8 C; R35	O; C R: 8-35 S: (1/2-)8-27-39-45		
011-004-00-7	sodium azide	247-852-1	26628-22-8	T+; R28 R32 N; R50-53	T+; N R: 28-32-50/53 S: (1/2-)28-45-60-61		
011-005-00-2	sodium carbonate	207-838-8	497-19-8	Xi; R36	Xi R: 36 S: (2-)22-26		
011-006-00-8	sodium cyanate	213-030-6	917-61-3	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)24/25-61		
011-007-00-3	propoxycarbazone-sodium	—	181274-15-7	N; R50-53	N R: 50/53 S: 60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	

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012-001-00-3	magnesium powder (pyrophoric)	231-104-6	7439-95-4	F; R15-17	F R: 15-17 S: (2-)7/8-43		
012-002-00-9	magnesium, powder or turnings	231-104-6	—	F; R11-15	F R: 11-15 S: (2-)7/8-43		
012-003-00-4	magnesium alkyls	—	—	R14 F; R17 C; R34	F; C R: 14-17-34 S: (1/2-)16-43-45		A
013-001-00-6	aluminium powder (pyrophoric)	231-072-3	7429-90-5	F; R15-17	F R: 15-17 S: (2-)7/8-43		
013-002-00-1	aluminium powder (stabilised)	231-072-3	—	F; R15 R10 ⊗	F R: 10-15 S: (2-)7/8-43		
013-003-00-7	aluminium chloride, anhydrous	231-208-1	7446-70-0	C; R34	C R: 34 S: (1/2-)7/8-28-45		
013-004-00-2	aluminium alkyls	—	—	R14 F; R17 C; R34	F; C R: 14-17-34 S: (1/2-)16-43-45		A
013-005-00-8	diethyl(ethyldimethylsilanolato)aluminium	401-160-8	55426-95-4	F; R 15-17 R14 C; R35	F; C R: 14/15-17-35 S: (1/2-)6-16-30-36/39-43-45		
013-006-00-3	(ethyl-3-oxobutanoato-O'1,O'3)(2-dimethylaminoethanolato)(1-methoxypropan-2-olato)aluminium(III), dimerised	402-370-2	—	R10 Xi; R41	Xi R: 10-41 S: (2-)26-39		
013-007-00-9	poly(oxo(2-butoxyethyl-3-oxobutanoato-O'1,O'3)aluminium)	403-430-0	—	Xi; R41	Xi R: 41 S: (2-)26-39		

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013-008-00-4	di-*n*-octylaluminium iodide	408-190-0	7585-14-0	R14 F; R17 C; R34 N; R50-53	F; C; N R: 14-17-34-50/53 S: (1/2-)6-16-26-36/37/39-43-45-60-61		
013-009-00-X	sodium((n-butyl)x(ethyl)y-1,5-dihydro)aluminate) x = 0,5, y = 1,5	418-720-2	—	F; R11-15-17 R14 Xn; R20 C; R35	F; C R: 11-14/15-17-20-35 S: (1/2-)6-16-26-30-36/37/39-43-45		
014-001-00-9	trichlorosilane	233-042-5	10025-78-2	F+; R12 R14 F; R17 Xn; R20/22 R29 C; R35	F+; C R: 12-14-17-20/22-29-35 S: (2-)7/9-16-26-36/37/39-43-45	Xn; R20/22: C ≥ 10 % C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
014-002-00-4	silicon tetrachloride	233-054-0	10026-04-7	R14 Xi; R36/37/38	Xi R: 14-36/37/38 S: (2-)7/8-26		
014-003-00-X	dimethyldichlorosilane	200-901-0	75-78-5	F; R11 Xi; R36/37/38	F; Xi R: 11-36/37/38 S: (2-)		
014-004-00-5	trichloro(methyl)silane; methyltrichlorosilane	200-902-6	75-79-6	R14 F; R11 Xi; R36/37/38	F; Xi R: 11-14-36/37/38 S: (2-)26-39	Xi; R36/37/38: C ≥ 1 %	
014-005-00-0	tetraethyl silicate; ethyl silicate	201-083-8	78-10-4	R10 Xn; R20 Xi; R36/37	Xn R: 10-20-36/37 S: (2-)		
014-006-00-6	bis(4-fluorophenyl)-methyl-(1,2,4-triazol-4-ylmethyl)silane hydrochloride	401-380-4	—	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
014-007-00-1	triethoxyisobutylsilane	402-810-3	17980-47-1	Xi; R38	Xi R: 38 S: (2-)24		

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014-008-00-7	(chloromethyl)bis(4-fluorophenyl)methylsilane	401-200-4	85491-26-5	N; R51-53	N R: 51/53 S: 61		
014-009-00-2	isobutylisopropyldimethoxysilane	402-580-4	111439-76-0	R10 Xn; R20 Xi; R38	Xn R: 10-20-38 S: (2-)25-26-36/37		
014-010-00-8	disodium metasilicate	229-912-9	6834-92-0	C; R34 Xi; R37	C R: 34-37 S: (1/2-)13-24/25-36/37/39-45		
014-011-00-3	cyclohexyldimethoxymethylsilane	402-140-1	17865-32-6	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)24-61		
014-012-00-9	bis(3-(trimethoxysilyl)propyl)amine	403-480-3	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)24-26-39-61		
014-013-00-4	α-hydroxypoly(methyl-(3-(2,2,6,6-tetramethylpiperidin-4-yloxy)propyl)siloxane)	404-920-7	—	Xn; R21/22 C; R34 N; R51-53	C; N R: 21/22-34-51/53 S: (1/2-)26-36/37/39-45-61		
014-014-00-X	etacelasil (ISO); 6-(2-chloroethyl)-6-(2-methoxyethoxy)-2,5,7,10-tetraoxa-6-silaundecane;	253-704-7	37894-46-5	Repr. Cat. 2; R61 Xn; R22-48/22	T R: 61-22-48/22 S: 53-45		E
014-015-00-5	α-trimethylsilanyl-ω-trimethylsiloxypoly[oxy(methyl-3-(2-(2-methoxypropoxy)propoxy)propylsilanediyl]-co-oxy(dimethyl-silane))	406-420-4	69430-40-6	R53	R: 53 S: 61		
014-016-00-0	reaction mass of: 1,3-dihex-5-en-1-yl-1,1,3,3-tetramethyldisiloxane; 1,3-dihex-n-en-1-yl-1,1,3,3-tetramethyldisiloxane	406-490-6	—	N; R51-53	N R: 51/53 S: 61		
014-017-00-6	flusilazole (ISO); bis(4-fluorophenyl)(methyl)(1*H*-1,2,4-triazol-1-ylmethyl) silane	—	85509-19-9	Carc. Cat. 3; R40 Repr. Cat. 2; R61 Xn; R22 N; R51-53	T; N R: 61-22-40-51/53 S: 53-45-61		E

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014-018-00-1	octamethylcyclotetrasiloxane	209-136-7	556-67-2	Repr. Cat. 3; R62 R53	Xn R: 53-62 S: (2-)36/37-46-51-61		
014-019-00-7	reaction mass of: 4-[[bis-(4-fluorophenyl)methylsilyl]methyl]-4*H*-1,2,4-triazole; 1-[[bis-(4-fluorophenyl)methylsilyl]methyl]-1*H*-1,2,4-triazole	403-250-2	—	Carc. Cat. 3; R40 Repr. Cat. 2; R61 Xn; R22 N; R51-53	T; N R: 61-22-40-51/53 S: 53-45-61		E
014-020-00-2	bis(1,1-dimethyl-2-propynyloxy)dimethylsilane	414-960-7	53863-99-3	Xn; R20	Xn R: 20 S: (2)		
014-021-00-8	tris(isopropenyloxy)phenyl silane	411-340-8	52301-18-5	N; R50-53	N R: 50/53 S: 60-61		
014-022-00-3	reaction product of: (2-hydroxy-4-(3-propenoxy)benzophenone and triethoxysilane) with (hydrolysis product of silica and methyltrimethoxysilane)	401-530-9	—	F; R11 T; R39/23/24/25 Xn; R20/21/22	F; T R: 11-20/21/22-39/23/24/25 S: (1/2-)16-29-36/37-45		
014-023-00-9	α, ω-dihydroxypoly(hex-5-en-1-ylmethylsiloxane)hoxysilane with (hydrolysis product of silica and methyltrimethoxysilane) iazole	408-160-7	125613-45-8	N; R51-53	N R: 51/53 S: 61		
014-024-00-4	1-((3-(3-chloro-4-fluorophenyl)propyl)dimethylsilanyl)-4-ethoxybenzene	412-620-2	121626-74-2	N; R51-53	N R: 51/53 S: 61		
014-025-00-X	4-[3-(diethoxymethylsilylpropoxy)-2,2,6,6-tetramethyl]piperidine	411-400-3	102089-33-8	Xn; R22-48/21 Xi; R38-41 R52-53	Xn R: 22-38-41-48/21-52/53 S: (2-)26-36/37/39-61		
014-026-00-5	dichloro-(3-(3-chloro-4-fluorophenyl)propyl)methylsilane	407-180-3	770722-36-6	C; R35	C R: 35 S: (1/2-)26-36/37/39-45		

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014-027-00-0	chloro(3-(3-chloro-4-fluorophenyl)propyl)dimethylsilane	410-270-5	770722-46-8	C; R35	C R: 35 S: (1/2-)8-26-28-36/37/39-45		
014-028-00-6	α-[3-(1-oxoprop-2-eny)l-1-oxypropyl]dimethoxysilyloxy-ω-[3(1-oxoprop-2-enyl)-1-oxypropyl]dimethoxysilyl poly(dimethylsiloxane)	415-290-8	193159-06-7	R43	Xi R: 43 S: (2-)24-37		
014-029-00-1	*O*,*O*'-(ethenylmethylsilylene)di[(4-methylpentan-2-one)oxime]	421-870-1	156145-66-3	Repr. Cat. 3; R62 Xn; R22-48/22	Xn R: 22-48/22-62 S: (2-)36/37		
014-030-00-7	[(dimethylsilylene)bis((1,2,3,3a,7a-η)-1*H*-inden-1-ylidene)dimethyl]hafnium	422-060-0	137390-08-0	T+; R28	T+ R: 28 S: (1/2-)6-22-28-36/37-45		
014-031-00-2	bis(1-methylethyl)-dimethoxysilane	421-540-7	18230-61-0	R10 Xi; R38 R43 R52-53	Xi R: 10-38-43-52/53 S: (2-)24-37-61		
014-032-00-8	dicyclopentyldimethoxysilane	404-370-8	126990-35-0	Xi; R38-41 N; R50-53	Xi; N R: 38-41-50/53 S: (2-)26-37/39-60-61		
015-001-00-1	white phosphorus	231-768-7	12185-10-3	F; R17 T+; R26/28 C; R35 N; R50	F; T+; C; N R: 17-26/28-35-50 S: (1/2-)5-26-38-45-61		
015-002-00-7	red phosphorus	231-768-7	7723-14-0	F; R11 R16 R52-53	F R: 11-16-52/53 S: (2-)7-43-61		
015-003-00-2	calcium phosphide; tricalcium diphosphide	215-142-0	1305-99-3	F; R15/29 T+; R28 N; R50	F; T+; N R: 15/29-28-50 S: (1/2-)22-43-45-61		

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015-004-00-8	aluminium phosphide	244-088-0	20859-73-8	F; R15/29 T+; R28 R32 N; R50	F; T+; N R: 15/29-28-32-50 S: (1/2-)3/9/14-30-36/37-45-61		
015-005-00-3	magnesium phosphide; trimagnesium diphosphide	235-023-7	12057-74-8	F; R15/29 T+; R28 N; R50	F; T+; N R: 15/29-28-50 S: (1/2-)22-43-45-61		
015-006-00-9	trizinc diphosphide; zinc phosphide	215-244-5	1314-84-7	F; R15/29 T+; R28 R32 N; R50-53	F; T+; N R: 15/29-28-32-50/53 S: (1/2-)3/9/14-30-36/37-45-60-61		
015-007-00-4	phosphorus trichloride	231-749-3	7719-12-2	R14 T+; R26/28 Xn; R48/20 C; R35 R29	T+; C R: 14-26/28-35-48/20 S: (1/2-)7/8-26-36/37/39-45		
015-008-00-X	phosphorus pentachloride	233-060-3	10026-13-8	R14 T+; R26 Xn; R22-48/20 C; R34 R29	T+ R: 14-22-26-34-48/20 S: (1/2-)7/8-26-36/37/39-45		
015-009-00-5	phosphoryl trichloride	233-046-7	10025-87-3	R14 T+; R26 T; R48/23 Xn; R22 C; R35 R29	T+; C R: 14-22-26-35-48/23 S: (1/2-)7/8-26-36/37/39-45		
015-010-00-0	phosphorus pentoxide	215-236-1	1314-56-3	C; R35	C R: 35 S: (1/2-)22-26-45		
015-011-00-6	phosphoric acid ... %, orthophosphoric acid ... %	231-633-2	7664-38-2	C; R34	C R: 34 S: (1/2-)26-45	C; R34: C ≥ 25 % Xi; R36/38: 10 % ≤ C < 25 %	B

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015-012-00-1	tetraphosphorus trisulphide; phosphorus sesquisulphid	215-245-0	1314-85-8	F; R11 Xn; R22 N; R50	F; Xn; N R: 11-22-50 S: (2-)7-16-24/25-61		
015-013-00-7	triethyl phosphate	201-114-5	78-40-0	Xn; R22	Xn R: 22 S: (2-)25		
015-014-00-2	tributyl phosphate	204-800-2	126-73-8	Carc. Cat. 3; R40 Xn; R22 Xi; R38	Xn R: 22-38-40 S: (2-)36/37-46		
015-015-00-8	tricresyl phosphate *(o-o-o-, o-o-m-, o-o-p-, o-m-m-, o-m-p-, o-p-p-)*; tritolyl phosphate *(o-o-o-, o-o-m-, o-o-p-, o-m-m-, o-m-p-, o-p-p-)*;	201-103-5	78-30-8	T; R39/23/24/25 N; R51-53	T; N R: 39/23/24/25-51/53 S: (1/2-)20/21-28-45-61	T; R39/23/24/25: C ≥ 1 % Xn; R68/20/21/22: 0,2 % ≤ C < 1 %	C
015-016-00-3	tricresyl phosphate *(m-m-m-, m-m-p-, m-p-p-, p-p-p-)*; tritolyl phosphate *(m-m-m-, m-m-p-, m-p-p-, p-p-p-)*;	201-105-6	78-32-0	Xn; R21/22 N; R51-53	Xn; N R: 21/22-51/53 S: (2-)28-61	Xn; R21/22: C ≥ 5 %	C
015-019-00-X	dichlorvos (ISO); 2,2-dichlorovinyl dimethyl phosphate	200-547-7	62-73-7	T+; R26 T; R24/25 R43 N; R50	T+; N R: 24/25-26-43-50 S: (1/2-)28-36/37-45-61		
015-020-00-5	mevinphos (ISO); 2-methoxycarbonyl-1-methylvinyl dimethyl phosphate	232-095-1	7786-34-7	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)23-28-36/37-45-60-61	N; R50-53: C ≥ 0,0025 % N; R51-53: 0,00025 % ≤ C < 0,0025 % R52-53: 0,000025 % ≤ C < 0,00025 %	
015-021-00-0	trichlorfon (ISO); dimethyl 2,2,2-trichloro-1-hydroxyethylphosphonate	200-149-3	52-68-6	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	

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015-022-00-6	phosphamidon (ISO); 2-chloro-2-diethylcarbamoyl-1-methylvinyl dimethyl phosphate	236-116-5	13171-21-6	T+; R28 T; R24 Muta. Cat. 3; R68 N; R50-53	T+; N R: 24-28-50/53-68 S: (1/2-)23-36/37-45-60-61		
015-023-00-1	pyrazoxon; diethyl 3-methylpyrazol-5-yl phosphate	—	108-34-9	T+; R26/27/28	T+ R: 26/27/28 S: (1/2-)13-28-45		
015-024-00-7	triamiphos (ISO); 5-amino-3-phenyl-1,2,4-triazol-1-yl-*N*,*N*,*N*',*N*'-tetramethylphosphonic diamide	—	1031-47-6	T+; R27/28	T+ R: 27/28 S: (1/2-)22-28-36/37-45		
015-025-00-2	TEPP (ISO); tetraethyl pyrophosphate	203-495-3	107-49-3	T+; R27/28 N; R50	T+; N R: 27/28-50 S: (1/2-)36/37/39-38-45-61		
015-026-00-8	schradan (ISO); octamethylpyrophosphoramide	205-801-0	152-16-9	T+; R27/28	T+ R: 27/28 S: (1/2-)36/37-38-45		
015-027-00-3	sulfotep (ISO); *O*,*O*,*O*,*O*-tetraethyl dithiopyrophosphate	222-995-2	3689-24-5	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)23-28-36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-028-00-9	demeton-*O* (ISO); *O*,*O*-diethyl-*O*-2-ethylthioethyl phosphorothioate	206-053-8	298-03-3	T+; R27/28 N; R50	T+; N R: 27/28-50 S: (1/2-)28-36/37-45-61		
015-029-00-4	demeton-*S* (ISO); diethyl-*S*-2-ethylthioethyl phosphorothioate	204-801-8	126-75-0	T+; R27/28	T+ R: 27/28 S: (1/2-)28-36/37-45		
015-030-00-X	demeton-*O*-methyl (ISO); *O*-2-ethylthioethyl *O*,*O*-dimethyl phosphorothioate	212-758-1	867-27-6	T; R25	T R: 25 S: (1/2-)24-36/37-45		

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015-031-00-5	demeton-*S*-methyl (ISO); *S*-2-ethylthioethyl dimethyl phosphorothioate	213-052-6	919-86-8	T; R24/25 N; R51-53	T; N R: 24/25-51/53 S: (1/2-)28-36/37-45-61		
015-032-00-0	prothoate (ISO); *O*,*O*-diethyl isopropylcarbamoylmethyl phosphorodithioate	218-893-2	2275-18-5	T+; R27/28 R52-53	T+ R: 27/28-52/53 S: (1/2-)28-36/37-45-61		
015-033-00-6	phorate (ISO); *O*,*O*-diethyl ethylthiomethyl phosphorodithioate	206-052-2	298-02-2	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-034-00-1	parathion (ISO); *O*,*O*-diethyl *O*-4-nitrophenyl phosphorothioate	200-271-7	56-38-2	T+; R26/28 T; R24-48/25 N; R50-53	T+; N R: 24-26/28-48/25-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-035-00-7	parathion — methyl (ISO); *O*,*O*-dimethyl *O*-4-nitrophenyl phosphorothioate	206-050-1	298-00-0	R5 R10 T+; R26/28 T; R24 Xn; R48/22 N; R50-53	T+; N R: 5-10-24-26/28-48/22-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-036-00-2	*O*-ethyl *O*-4-nitrophenyl phenylphosphonothioate; EPN	218-276-8	2104-64-5	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)22-36/37-45-60-61		
015-037-00-8	phenkapton (ISO); *S*-(2,5-dichlorophenylthiomethyl) *O*,*O*-diethyl phosphorodithioate	218-892-7	2275-14-1	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		
015-038-00-3	coumaphos (ISO); *O*-3-chloro-4-methylcoumarin-7-yl *O*,*O*-diethyl phosphorothioate	200-285-3	56-72-4	T+; R28 Xn; R21 N; R50-53	T+; N R: 21-28-50/53 S: (1/2-)28-36/37-45-60-61		

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015-039-00-9	azinphos-methyl (ISO); *O*,*O*-dimethyl-4-oxobenzotriazin-3-ylmethyl phosphorodithioate	201-676-1	86-50-0	T+; R26/28 T; R24 R43 N; R50-53	T+; N R: 24-26/28-43-50/53 S: (1/2-)28-36/37-45-60-61		
015-040-00-4	diazinon (ISO); *O*,*O*-diethyl *O*-2-isopropyl-6-methylpyrimidin-4-yl phosphorothioate	206-373-8	333-41-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24/25-60-61		
015-041-00-X	malathion (ISO); 1,2-bis (ethoxycarbonyl) ethyl *O*,*O*-dimethyl phosphorodithioate	204-497-7	121-75-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-042-00-5	chlorthion *O*-(3-chloro-4-nitrophenyl) *O*,*O*-dimethyl phosphorothioate	207-902-5	500-28-7	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-043-00-0	phosnichlor (ISO); *O*-4-chloro-3-nitrophenyl *O*,*O*-dimethyl phosphorothioate	—	5826-76-6	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)13		
015-044-00-6	carbophenothion (ISO); 4-chlorophenylthiomethyl *O*,*O*-diethyl phosphorodithioate	212-324-1	786-19-6	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61		
015-045-00-1	mecarbam (ISO); *N*-ethoxycarbonyl-*N*-methylcarbamoylmethyl *O*,*O*-diethyl phosphorodithioate	219-993-9	2595-54-2	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)36/37-45-60-61		
015-046-00-7	oxydemeton-methyl; *S*-2-(ethylsulphinyl)ethyl *O*,*O*-dimethyl phosphorothioate	206-110-7	301-12-2	T; R24/25 N; R50	T; N R: 24/25-50 S: (1/2-)23-36/37-45-61		
015-047-00-2	ethion (ISO); *O*,*O*,*O*',*O*'-tetraethyl *S*,*S*'-methylenedi (phosphorodithioate); diethion	209-242-3	563-12-2	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)25-36/37-45-60-61	N; R50-53: C ≥ 0,0025 % N; R51-53: 0,00025 % ≤ C < 0,0025 % R52-53: 0,000025 % ≤ C < 0,00025 %	

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015-048-00-8	fenthion (ISO); *O*,*O*-dimethyl-*O*-(4-methylthion-*m*-tolyl) phosphorothioate	200-231-9	55-38-9	Muta. Cat. 3; R68 T; R23-48/25 Xn; R21/22 N; R50-53	T; N R: 21/22-23-48/25-50/53-68 S: (1/2-)36/37-45-60-61		
015-049-00-3	endothion (ISO); *S*-5-methoxy-4-oxopyran-2-ylmethyl dimethyl phosphorothioate	220-472-3	2778-04-3	T; R24/25	T R: 24/25 S: (1/2-)36/37-45		
015-050-00-9	thiometon (ISO); *S*-2-ethylthioethyl *O*,*O*-dimethyl phosphorodithioate	211-362-6	640-15-3	T; R25 Xn; R21	T R: 21-25 S: (1/2-)36/37-45		
015-051-00-4	dimethoate (ISO); *O*,*O*-dimethyl methylcarbamoylmethyl phosphorodithioate	200-480-3	60-51-5	Xn; R21/22	Xn R: 21/22 S: (2-)36/37		
015-052-00-X	fenchlorphos (ISO); *O*,*O*-dimethyl *O*-2,4,5-trichlorophenyl phosphorothioate	206-082-6	299-84-3	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)25-36/37-60-61		
015-053-00-5	menazon (ISO); S-[(4,6-diamino-1,3,5-triazin-2-yl)methyl] O,O-dimethyl phosphorodithioate	201-123-4	78-57-9	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
015-054-00-0	fenitrothion (ISO); *O*,*O*-dimethyl *O*-4-nitro-*m*-tolyl phosphorothioate	204-524-2	122-14-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
015-055-00-6	naled (ISO); 1,2-dibromo-2,2-dichloroethyl dimethyl phosphate	206-098-3	300-76-5	Xn; R21/22 Xi; R36/38 N; R50	Xn; N R: 21/22-36/38-50 S: (2-)36/37-61	N; R50: C ≥ 0,025 %	
015-056-00-1	azinphos-ethyl (ISO); *O*,*O*-diethyl 4-oxobenzotriazin-3-ylmethyl phosphorodithioate	220-147-6	2642-71-9	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61		
015-057-00-7	formothion (ISO); *N*-formyl-*N*-methylcarbamoylmethyl *O*,*O*-dimethyl phosphorodithioate	219-818-6	2540-82-1	Xn; R21/22	Xn R: 21/22 S: (2-)36/37		

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015-058-00-2	morphothion (ISO); *O,O*-dimethyl-*S*-(morpholinocarbonylmethyl) phosphorodithioate	205-628-0	144-41-2	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		
015-059-00-8	vamidothion (ISO); *O,O*-dimethyl *S*-2-(1-methylcarbamoylethylthio) ethyl phosphorothioate	218-894-8	2275-23-2	T; R25 Xn; R21 N; R50	T; N R: 21-25-50 S: (1/2-)36/37-45-61		
015-060-00-3	disulfoton (ISO); *O,O*-diethyl 2-ethylthioethyl phosphorodithioate	206-054-3	298-04-4	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)28-36/37-45-60-61		
015-061-00-9	dimefox (ISO); tetramethylphosphorodiamidic fluoride	204-076-8	115-26-4	T+; R27/28	T+ R: 27/28 S: (1/2-)23-28-36/37-38-45		
015-062-00-4	mipafox (ISO); *N,N'*- di-isopropylphosphorodiamidic fluoride	206-742-3	371-86-8	T+; R39/26/27/28	T+ R: 39/26/27/28 S: (1/2-)13-45		
015-063-00-X	dioxathion (ISO); 1,4-dioxan-2,3-diyl-*O,O,O',O'*-tetraethyl di(phosphorodithioate)	201-107-7	78-34-2	T+; R26/28 T; R24 N; R50-53	T+; N R: 24-26/28-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-064-00-5	bromophos-ethyl (ISO); *O*-4-bromo-2,5-dichlorophenyl *O,O*-diethyl phosphorothioate	225-399-0	4824-78-6	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)28-36/37-45-60-61		
015-065-00-0	*S*-[2-(ethylsulphinyl)ethyl] *O,O*-dimethyl phosphorodithioate	—	2703-37-9	T+; R26/27/28 N; R51-53	T+; N R: 26/27/28-51/53 S: (1/2-)13-28-45-61		
015-066-00-6	omethoate (ISO); *O,O*-dimethyl *S*-methylcarbamoylmethyl phosphorothioate	214-197-8	1113-02-6	T; R25 Xn; R21 N; R50	T; N R: 21-25-50 S: (1/2-)23-36/37-45-61		

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015-067-00-1	phosalone (ISO); *S*-(6-chloro-2-oxobenzoxazolin-3-ylmethyl) *O*,*O*-diethyl phosphorodithioate	218-996-2	2310-17-0	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)36/37-45-60-61		
015-068-00-7	dichlofenthion (ISO); *O*-,4-dichlorophenyl *O*,*O*-diethyl phosphorothioate	202-564-5	97-17-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
015-069-00-2	methidathion (ISO); 2,3-dihydro-5-methoxy-2-oxo-1,3,4-thiadiazol-3-ylmethyl-*O*, *O*-dimethylphosphorodithioate	213-449-4	950-37-8	T+; R28 Xn; R21 N; R50-53	T+; N R: 21-28-50/53 S: (1/2-)22-28-36/37-45-60-61		
015-070-00-8	cyanthoate (ISO); *S*-(*N*-(1-cyano-1-methylethyl)carbamoylmethyl) *O*,*O*-diethyl phosphorothioate	223-099-4	3734-95-0	T+; R28 T; R24	T+ R: 24-28 S: (1/2-)36/37-45		
015-071-00-3	chlorfenvinphos (ISO); 2-chloro-1-(2,4 dichlorophenyl) vinyl diethyl phosphate	207-432-0	470-90-6	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61		
015-072-00-9	monocrotophos (ISO); dimethyl-1-methyl-2-(methylcarbamoyl)vinyl phosphate	230-042-7	6923-22-4	Muta. Cat. 3; R68 T+; R26/28 T; R24 N; R50-53	T+; N R: 24-26/28-50/53-68 S: (1/2-)36/37-45-60-61		
015-073-00-4	dicrotophos (ISO); (*Z*)-2-dimethylcarbamoyl-1-methylvinyl dimethyl phosphate	205-494-3	141-66-2	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61		
015-074-00-X	crufomate (ISO); 4-tert-butyl-2-chlorophenyl methyl methylphosphoramidate	206-083-1	299-86-5	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
015-075-00-5	*S*-[2-(isopropylsulphinyl)ethyl] *O*,*O*-dimethyl phosphorothioate	—	2635-50-9	T; R23/24/25	T R: 23/24/25 S: (1/2-)13-45		

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015-076-00-0	potasan; *O*,*O*-diethyl *O*-(4-methylcoumarin-7-yl) phosphorothioate	—	299-45-6	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)13-28-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-077-00-6	2,2-dichlorovinyl 2-ethylsulphinylethyl methyl phosphate	—	7076-53-1	T; R23/24/25	T R: 23/24/25 S: (1/2-)13-45		
015-078-00-1	demeton-*S*-methylsulphon (ISO); *S*-2-ethylsulphonylethyl dimethyl phosphorothioate	241-109-5	17040-19-6	T; R25 Xn; R21 N; R51-53	T; N R: 21-25-51/53 S: (1/2-)22-28-36/37-45-61		
015-079-00-7	acephate (ISO); *O*,*S*-dimethyl acetylphosphoramidothioate	250-241-2	30560-19-1	Xn; R22	Xn R: 22 S: (2-)36		
015-080-00-2	amidithion (ISO); 2-methoxyethylcarbamoylmethyl *O*,*O*-dimethyl phosphorodithioate	—	919-76-6	Xn; R22	Xn R: 22 S: (2-)24-36		
015-081-00-8	*O*,*O*,*O*',*O*'-tetrapropyl dithiopyrophosphate	221-817-0	3244-90-4	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
015-082-00-3	azothoate (ISO); *O*-4-(4-chlorophenylazo)phenyl *O*,*O*-dimethyl phosphorothioate	227-419-3	5834-96-8	Xn; R20/22	Xn R: 20/22 S: (2-)13		
015-083-00-9	bensulide (ISO); *O*,*O*-diisopropyl 2-phenylsulphonylaminoethyl phosphorodithioate	212-010-4	741-58-2	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24-36-60-61		
015-084-00-4	chlorpyrifos (ISO); *O*,*O*-diethyl *O*-3,5,6-trichloro-2-pyridyl phosphorothioate	220-864-4	2921-88-2	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)45-60-61	N; R50-53: C ≥ 0,0025 % N; R51-53: 0,00025 % ≤ C < 0,0025 % R52-53: 0,000025 % ≤ C < 0,00025 %	
015-085-00-X	chlorphonium chloride (ISO); tributyl (2,4-dichlorobenzyl) phosphonium chloride	204-105-4	115-78-6	T; R25 Xn; R21 Xi; R36/38	T R: 21-25-36/38 S: (1/2-)36/37/39-45		

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015-086-00-5	coumithoate (ISO); *O*,*O*-diethyl *O*-,8,9,10-tetrahydro-6-oxo-benzo(c)chromen-3-yl phosphorothioate	—	572-48-5	T; R25	T R: 25 S: (1/2-)28-36/37-45		
015-087-00-0	cyanophos (ISO); *O*-4-cyanophenyl *O*,*O*-dimethyl phosphorothioate	220-130-3	2636-26-2	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
015-088-00-6	dialifos (ISO); 2-chloro-1-phthalimidoethyl *O*,*O*-diethyl phosphorodithioate	233-689-3	10311-84-9	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61		
015-089-00-1	ethoate-methyl (ISO); ethylcarbamoylmethyl *O*,*O*-dimethyl phosphorodithioate	204-121-1	116-01-8	Xn; R21/22	Xn R: 21/22 S: (2-)36/37		
015-090-00-7	fensulfothion (ISO); *O*,*O*-diethyl *O*-4-methylsulfinylphenyl phosphorothioate	204-114-3	115-90-2	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)23-28-36/37-45-60-61		
015-091-00-2	fonofos (ISO); *O*-ethyl phenyl ethylphosphonodithioate	213-408-0	944-22-9	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)28-36/37-45-60-61		
015-092-00-8	phosacetim (ISO); *O*,*O*-bis(4-chlorophenyl) *N*-acetimidoylphosphoramidothioate	223-874-7	4104-14-7	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)28-36/37-45-60-61		
015-093-00-3	leptophos (ISO); *O*-4-bromo-2,5-dichlorophenyl *O*-methyl phenylphosphorothioate	244-472-8	21609-90-5	T; R25-39/25 Xn; R21 N; R50-53	T; N R: 21-25-39/25-50/53 S: (1/2-)25-36/37/39-45-60-61		
015-094-00-9	mephosfolan (ISO); diethyl 4-methyl-1,3-dithiolan-2-ylidenephosphoramidate	213-447-3	950-10-7	T+; R27/28 N; R51-53	T+; N R: 27/28-51/53 S: (1/2-)36/37/39-45-61		

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015-095-00-4	methamidophos (ISO); *O,S*-dimethyl phosphoramidothioate	233-606-0	10265-92-6	T+; R26/28 T; R24 N; R50	T+; N R: 24-26/28-50 S: (1/2-)28-36/37-45-61		
015-096-00-X	oxydisulfoton (ISO); *O, O*-diethyl *S*-2-ethylsulphinylethyl phosphorodithioate	219-679-1	2497-07-6	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
015-097-00-5	phenthoate (ISO); ethyl 2-(dimethoxyphosphinothioylthio)-2-phenylacetate	219-997-0	2597-03-7	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)22-36/37-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-098-00-0	trichloronate (ISO); *O*-ethyl *O*-2,4,5-trichlorophenyl ethylphosphonothioate	206-326-1	327-98-0	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)23-28-36/37-45-60-61		
015-099-00-6	pirimiphos-ethyl (ISO); *O,O*-diethyl *O*-2-diethylamino-6-methylpyrimidin-4-yl phosphorothioate	245-704-0	23505-41-1	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)23-36/37-45-60-61		
015-100-00-X	phoxim (ISO); α-(diethoxyphosphinothioylimino) phenylacetonitrile	238-887-3	14816-18-3	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)36-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-101-00-5	phosmet (ISO); *O,O*-dimethyl phthalimidomethyl *S*-phosphorodithioate	211-987-4	732-11-6	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)22-36/37-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-102-00-0	tris(2-chloroethyl) phosphate	204-118-5	115-96-8	Carc. Cat. 3; R40 Xn; R22 N; R51-53	Xn; N R: 22-40-51/53 S: (2-)36/37-61		
015-103-00-6	phosphorus tribromide	232-178-2	7789-60-8	R14 C; R34 Xi; R37	C R: 14-34-37 S: (1/2-)26-45		

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015-104-00-1	diphosphorus pentasulphide; phosphorus pentasulphide	215-242-4	1314-80-3	F; R11 R29 Xn; R20/22 N; R50	F; Xn; N R: 11-20/22-29-50 S: (2-)61		
015-105-00-7	triphenyl phosphite	202-908-4	101-02-0	Xi; R36/38 N; R50-53	Xi; N R: 36/38-50/53 S: (2-)28-60-61	Xi; R36/38: C ≥ 5 %	
015-106-00-2	hexamethylphosphoric triamide; hexamethylphosphoramide	211-653-8	680-31-9	Carc. Cat. 2; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45	Carc. Cat. 2; R45: C ≥ 0,01 %	
015-107-00-8	ethoprophos (ISO); ethyl-*S*,*S*-dipropyl phosphorodithioate	236-152-1	13194-48-4	T+; R26/27 T; R25 R43 N; R50-53	T+; N R: 25-26/27-43-50/53 S: (1/2-)27/28-36/37/39-45-60-61		
015-108-00-3	bromophos (ISO); *O*-4-bromo-2,5-dichlorophenyl *O*,*O*-dimethyl phosphorothioate	218-277-3	2104-96-3	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-) 46-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-109-00-9	crotoxyphos (ISO); 1-phenylethyl 3-(dimethoxyphosphinyloxy) isocrotonate	231-720-5	7700-17-6	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
015-110-00-4	cyanofenphos (ISO); *O*-4-cyanophenyl *O*-ethyl phenylphosphonothioate	—	13067-93-1	T; R25-39/25 Xn; R21 Xi; R36 N; R51-53	T; N R: 21-25-36-39/25-51/53 S: (1/2-)36/37-45-61		
015-111-00-X	phosfolan (ISO); diethyl 1,3-dithiolan-2-ylidenephosphoramidate	213-423-2	947-02-4	T+; R27/28	T+ R: 27/28 S: (1/2-)28-36/37-45		
015-112-00-5	thionazin (ISO); *O*,*O*-diethyl *O*-pyrazin-2-yl phosphorothioate;	206-049-6	297-97-2	T+; R27/28	T+ R: 27/28 S: (1/2-)36/37/39-38-45		

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015-114-00-6	chlormephos (ISO); *S*-chloromethyl *O*,*O*-diethyl phosphorodithioate	246-538-1	24934-91-6	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)28-36/37-45-60-61		
015-115-00-1	chlorthiophos (ISO)	244-663-6	21923-23-9	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)28-36/37-45-60-61		
015-116-00-7	demephion-*O* (ISO); *O*,*O*-dimethyl *O*-2-methylthioethyl phosphorothioate	211-666-9	682-80-4	T+; R28 T; R24	T+ R: 24-28 S: (1/2-)28-36/37-45		
015-117-00-2	demephion-*S* (ISO); *O*,*O*-dimethyl *S*-2-methylthioethyl phosphorothioate	219-971-9	2587-90-8	T+; R28 T; R24	T+ R: 24-28 S: (1/2-)28-36/37-45		
015-118-00-8	demeton	—	8065-48-3	T+; R27/28 N; R50	T+; N R: 27/28-50 S: (1/2-)28-36/37-45-61		
015-119-00-3	dimethyl 4-(methylthio)phenyl phosphate	—	3254-63-5	T+; R27/28	T+ R: 27/28 S: (1/2-)28-36/37-45		
015-120-00-9	ditalimfos (ISO); *O*,*O*-diethyl phthalimidophosphonothioate;	225-875-8	5131-24-8	Xi; R38 R43	Xi R: 38-43 S: (2-)36/37		
015-121-00-4	edifenphos (ISO); *O*-ethyl *S*,*S*-diphenyl phosphorodithioate	241-178-1	17109-49-8	T; R23/25 Xn; R21 R43 N; R50-53	T; N R: 21-23/25-43-50/53 S: (1/2-)36/37-45-60-61		
015-122-00-X	etrimfos (ISO); *O*-6-ethoxy-2-ethylpyrimidin-4-yl *O*,*O*-dimethylphosphorothioate;	253-855-9	38260-54-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	

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015-123-00-5	fenamiphos (ISO); ethyl-4-methylthio-*m*-tolyl isopropyl phosphoramidate	244-848-1	22224-92-6	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)23-28-36/37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-124-00-0	fosthietan (ISO); diethyl 1,3-dithietan-2-ylidenephosphoramidate;	244-437-7	21548-32-3	T+; R27/28	T+ R: 27/28 S: (1/2-)36/37-45		
015-125-00-6	glyphosine (ISO); *N*,*N*-bis(phosphonomethyl)glycine	219-468-4	2439-99-8	Xi; R41	Xi R: 41 S: (2-)26		
015-126-00-1	heptenophos (ISO); 7-chlorobicyclo(3.2.0)hepta-2,6-dien-6-yl dimethyl phosphate	245-737-0	23560-59-0	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)23-28-37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-127-00-7	iprobenfos (ISO); *S*-benzyl diisopropyl phosphorothioate	247-449-0	26087-47-8	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
015-128-00-2	IPSP; *S*-ethylsulphinylmethyl *O*,*O*-diisopropylphosphorodithioate	—	5827-05-4	T+; R27 T; R25 N; R50-53	T+; N R: 25-27-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-129-00-8	isofenphos (ISO); *O*-ethyl *O*-2-isopropoxycarbonylphenyl-isopropylphosphoramidothioate	246-814-1	25311-71-1	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)36/37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-130-00-3	isothioate (ISO) *S*-2-isopropylthioethyl *O*,*O*-dimethyl phosphorodithioate;	—	36614-38-7	T; R24/25	T R: 24/25 S: (1/2-)28-36/37-45		
015-131-00-9	isoxathion (ISO); *O*,*O*-diethyl *O*-5-phenylisoxazol-3-ylphosphorothioate	242-624-8	18854-01-8	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61		

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015-132-00-4	*S*-(chlorophenylthiomethyl) *O*,*O*-dimethylphosphorodithioate; methylcarbophenothione	—	953-17-3	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)28-36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-133-00-X	piperophos (ISO); *S*-2-methylpiperidinocarbonylmethyl-*O*,*O*-dipropyl phosphorodithioate	—	24151-93-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
015-134-00-5	pirimiphos-methyl (ISO); *O*-(2-diethylamino-6-methylpyrimidin-4-yl) *O*,*O*-dimethyl phosphorothioate	249-528-5	29232-93-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
015-135-00-0	profenofos (ISO); *O*-(4-bromo-2-chlorophenyl) *O*-ethyl *S*-propyl phosphorothioate	255-255-2	41198-08-7	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)36/37-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-136-00-6	*trans*-isopropyl-3-[[(ethylamino)methoxyfosfinothioyl]oxy] crotonate; isopropyl 3-[[(ethylamino)methoxyphosphinothioyl]oxy]isocrotonate; propetamphos (ISO)	250-517-2	31218-83-4	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)37-45-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
015-137-00-1	pyrazophos (ISO); *O*,*O*-diethyl *O*-(6-ethoxycarbonyl-5-methylpyrazolo[2,3-a] pyrimidin-2-yl) phosphorothioate	236-656-1	13457-18-6	Xn; R20/22 N; R50-53	Xn; N R: 20/22-50/53 S: (2-)36/37-46-60-61		
015-138-00-7	quinalphos (ISO); *O*,*O*-diethyl-*O*-quinoxalin-2-yl phosphorothioate	237-031-6	13593-03-8	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)22-36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-139-00-2	terbufos (ISO) *S*-*tert*-butylthiomethyl *O*,*O*-diethylphosphorodithioate;	235-963-8	13071-79-9	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)36/37-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
015-140-00-8	triazophos (ISO); *O*,*O*-diethyl-*O*-1-phenyl-1*H*,2,4-triazol-3-yl phosphorothioate	245-986-5	24017-47-8	T; R23/25 Xn; R21 N; R50-53	T; N R: 21-23/25-50/53 S: (1/2-)36/37-45-60-61		

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015-141-00-3	ethylenediammonium *O,O*-bis(octyl) phosphorodithioate, mixed isomers	400-520-1	—	C; R34 Xn; R22 N; R50-53	C; N R: 22-34-50/53 S: (1/2-)24/25-26-28-39-45-60-61		
015-142-00-9	butyl (dialkyloxy(dibutoxyphosphoryloxy))titanium (trialkyloxy)titanium phosphate	401-100-0	—	F; R11 Xi; R36 N; R51-53	F; Xi; N R: 11-36-51/53 S: (2-)7/9-16-26-43-61		
015-143-00-4	reaction mass of 2-chloroethyl chloropropyl 2-chloroethylphosphonate, ~~mixture~~ reaction mass of isomers and 2-chloroethyl chloropropyl 2-chloropropylphosphonate, reaction mass of isomers	401-740-0	—	Xn; R22	Xn R: 22 S: (2-)		
015-144-00-X	reaction mass of pentyl methylphosphinate and 2-methylbutyl methylphosphinate	402-090-0	87025-52-3	C; R34	C R: 34 S: (1/2-)26-36/37/39-45		
015-145-00-5	reaction mass of copper(I) *O,O*-diisopropyl phosphorodithioate and copper(I) *O*-isopropyl *O*-(4-methylpent-2-yl) phosphorodithioate and copper(I) *O,O*-bis(4-methylpent-2-yl) phosphorodithioate	401-520-4	—	N; R50-53	N R: 50/53 S: 60-61		
015-146-00-0	*S*-(tricyclo(5.2.1.$0^{2,6}$)deca-3-en-8(or 9)-yl *O*-(isopropyl or isobutyl or 2-ethylhexyl) *O*-(isopropyl or isobutyl or 2-ethylhexyl) phosphorodithioate	401-850-9	—	N; R50-53	N R: 50/53 S: 60-61		
015-147-00-6	reaction mass of C_{12-14}-tert-alkylammonium diphenyl phosphorothioate and dinonyl sulphide (or disulphide)	400-930-0	—	Xi; R38-41 N; R51-53 R43	Xi; N R: 38-41-43-51/53 S: (2-)24-26-37/39-61		
015-148-00-1	2-(diphosphonomethyl)succinic acid	403-070-4	51395-42-7	C; R34 R43	C R: 34-43 S: (1/2-)26-36/37/39-45		
015-149-00-7	reaction mass of: hexyldioctylphosphineoxide; dihexyloctylphosphineoxide; trioctylphosphineoxide	403-470-9	—	C; R34 N; R50-53	C; N R: 34-50/53 S: (1/2-)26-36/37/39-45-60-61		

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015-150-00-2	(2-(1,3-dioxolan-2-yl)ethyl)triphenylphosphonium bromide	404-940-6	86608-70-0	Xn; R22 Xi; R41 R33 R52-53	Xn R: 22-33-41-52/53 S: (2-)22-26-39-61		
015-151-00-8	tris(isopropyl/*tert*-butylphenyl) phosphate	405-010-2	—	N; R51-53	N R: 51/53 S: 61		
015-152-00-3	dioxabenzofos (ISO); 2-methoxy-4*H*-1,3,2-benzodioxaphosphorin 2-sulphide;	223-292-3	3811-49-2	T; R24/25-39/25 N; R51-53	T; N R: 24/25-39/25-51/53 S: (1/2-)36/37-38-45-61		
015-153-00-9	isazofos (ISO); *O*-(5-chloro-1-isopropyl-1,2,4-triazol-3-yl) *O*,*O*-diethyl phosphorothioate;	255-863-8	42509-80-8	T+; R26 T; R24/25 Xn; R48/20 R43 N; R50-53	T+; N R: 24/25-26-43-48/20-50/53 S: (1/2-)28-36/37-38-45-59-61		
015-154-00-4	2-chloroethylphosphonic acid; ethephon	240-718-3	16672-87-0	Xn; R20/21 C; R34 R52-53	C R: 20/21-34-52/53 S: (1/2-)26-28-36/37/39-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
015-155-00-X	ammonium 2-amino-4-(hydroxymethylphosphinyl)butyrate; glufosinate ammonium	278-636-5	77182-82-2	Xn; R22	Xn R: 22 S: (2-)		
015-156-00-5	methyl 3-[(dimethoxyphosphinothioyl)oxy]methacrylate; [1] methacrifos (ISO); methyl (*E*)-3-[(dimethoxyphosphinothioyl)oxy]methacrylate [2]	250-366-9 [1] - [2]	30864-28-9 [1] 62610-77-9 [2]	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)36/37-60-61		
015-157-00-0	phosphonic acid; [1] phosphorous acid [2]	237-066-7 [1] 233-663-1 [2]	13598-36-2 [1] 10294-56-1 [2]	Xn; R22 C; R35	C R: 22-35 S: (1/2-)26-36/37/39-45		
015-158-00-6	(η-cyclopentadienyl)(η-cumenyl)iron(1+)hexafluorophosphate (1-)	402-340-9	32760-80-8	R52-53	R: 52/53 S: 61		

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015-159-00-1	hydroxyphosphonoacetic acid	405-710-8	23783-26-8	Xn; R22-48/22 C; R34 R43	C R: 22-34-43-48/22 S: (1/2-)22-26-36/37/39-45		
015-160-00-7	vanadyl pyrophosphate	406-260-5	58834-75-6	Xi; R36 R43 R52-53	Xi R: 36-43-52/53 S: (2-)24-26-37-61		
015-161-00-2	divanadyl pyrophosphate	407-130-0	65232-89-5	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)24-26-37/39-61		
015-162-00-8	vanadium(IV) oxide hydrogen phosphate hemihydrate, lithium, zinc, molybdenum, iron and chlorine-doped	407-350-7	—	Xn; R20-48/22 Xi; R41 N; R51-53	Xn; N R: 20-41-48/22-51/53 S: (2-)22-26-36/39-61		
015-163-00-3	bis(2,6-dimethoxybenzoyl)-2,4,4-trimethylpentylphosphinoxide	412-010-6	145052-34-2	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
015-164-00-9	calcium *P,P'*-(1-hydroxyethylene)bis(hydrogen phosphonate) dihydrate	400-480-5	36669-85-9	R52-53	R: 52/53 S: 61		
015-165-00-4	reaction mass of: thiobis(4,1-phenylene)-*S,S,S',S'*-tetraphenyldisulfonium bishexafluorophosphate; diphenyl(4-phenylthiophenyl)sulfonium hexafluorophosphate	404-986-7	—	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)15-26-39-60-61		
015-166-00-X	3,9-bis(2,6-di-*tert*-butyl-4-methylphenoxy)-2,4,8,10-tetraoxa-3,9-diphosphaspiro[5.5]undecane	410-290-4	80693-00-1	R53	R: 53 S: 61		
015-167-00-5	3-(hydroxyphenylphosphinyl)propanoic acid	411-200-6	14657-64-8	Xi; R41	Xi R: 41 S: (2-)26-39		
015-168-00-0	fosthiazate (ISO); (*RS*)-*S*-*sec*-butyl-O-ethyl-2-oxo-1,3-thiazolidin-3-ylphosphonothioate	—	98886-44-3	T; R23/25-39 Xn; R21 Xi; R41 R43 N; R50-53	T; N R: 21-23/25-39-41-43-50/53 S: (1/2-)53-45-25-26-39-60-61		

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015-169-00-6	tributyltetradecylphosphonium tetrafluoroborate	413-520-1	—	Xn; R22-48/22 C; R34 R43 N; R50-53	C; N R: 22-34-43-48/22-50/53 S: (1/2-)26-28-36/37/39-45-60-61		
015-170-00-1	reaction mass of: di-(1-octane-*N*,*N*,*N*-trimethylammonium) octylphosphate; 1-octane-*N*,*N*,*N*-trimethylammonium di-octylphosphate; 1-octane-*N*,*N*,*N*-trimethylammonium octylphosphate	407-490-9	—	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-36/37/39-45		
015-171-00-7	*O*,*O*,*O*-tris(2(or 4)-C_{9-10}-isoalkylphenyl) phosphorothioate	406-940-1	—	N; R51-53	N R: 51/53 S: 61		
015-172-00-2	reaction mass of: bis(isotridecylammonium)mono (di-(4-methylpent-2-yloxy)thiophosphorothionylisopropyl) phosphate; isotridecylammonium bis(di-(4-methylpent-2-yloxy) thiophosphorothionylisopropyl)phosphate	406-240-6	—	R10 C; R34 N; R51-53	C; N R: 10-34-51/53 S: (1/2-)23-26-28-36/37/39-45-61		
015-173-00-8	methyl [2-(1,1-dimethylethyl)-6-methoxypyrimidin-4-yl] ethylphosphonothioate	414-080-3	117291-73-3	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)23-36-60-61		
015-174-00-3	1-chloro-*N*,*N*-diethyl-1,1-diphenyl-1-(phenylmethyl)phosphoramine	411-370-1	82857-68-9	T; R25 Xi; R41 N; R51-53	T; N R: 25-41-51/53 S: (1/2-)26-37/39-41-45-61		
015-175-00-9	*tert*-butyl (triphenylphosphoranylidene) acetate	412-880-7	35000-38-5	T; R25 Xn; R48/22 Xi; R36 R43 N; R51-53	T; N R: 25-36-43-48/22-51/53 S: (1/2-)26-36/37/39-45-61		
015-176-00-4	*P*,*P*,*P*′,*P*′-tetrakis-(*o*-methoxyphenyl)propane-1,3-diphosphine	413-430-2	116163-96-3	N; R50-53	N R: 50/53 S: 60-61		
015-177-00-X	((4-phenylbutyl)hydroxyphosphoryl)acetic acid	412-170-7	83623-61-4	Xn; R48/22 Xi; R41 R43	Xn R: 41-43-48/22 S: (2-)22-26-36/37/39		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
015-178-00-5	(*R*)-α-phenylethylammonium (-)-(1*R*, 2*S*)-(1,2-epoxypropyl) phosphonate monohydrate	418-570-8	25383-07-7	Repr. Cat. 3; R62 N; R51-53	Xn; N R: 62-51/53 S: (2-)22-36/37-61		
015-179-00-0	UVCB condensation product of: tetrakis-hydroxymethyl-phosphonium chloride, urea and distilled hydrogenated C_{16-18} tallow alkylamine	422-720-8	166242-53-1	Carc. Cat. 3; R40 Xn; R22-48/22 C; R34 R43 N; R50-53	C; N R: 22-34-40-43-48/22-50/53 S: (1/2-)26-36/37/39-45-60-61		
015-180-00-6	[*R*-(*R**,*S**)]-[[2-methyl-1-(1-oxopropoxy)propoxy]-(4-phenylbutyl)phosphinyl] acetic acid, (-)-cinchonidine (1:1) salt	415-820-8	137590-32-0	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
015-181-00-1	phosphine	232-260-8	7803-51-2	F+; R12 R17 T+; R26 C; R34 N; R50	F+; T+; N R: 12-17-26-34-50 S: (1/2-)28-36/37-45-61-63		
015-184-00-8	Salts of glyphosate, with the exception of those specified elsewhere in this Annex	—	—	N; R51-53	N R: 51/53 S: 61		A
015-186-00-9	chlorpyrifos-methyl (ISO), *O*, *O*-dimethyl *O*-3,5,6-trichloro-2-pyridyl phosphorothioate	227-011-5	5598-13-0	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)36/37-60-61	N; R50-53: C ≥ 0,0025 % N; R51-53: 0,00025 % ≤ C < 0,0025 % R52-53: 0,000025 % ≤ C < 0,00025 %	
015-187-00-4	reaction mass of: tetrasodium(((2-hydroxyethyl)imino)bis (methylene))bisphosphonate, *N*-oxide; trisodium ((tetrahydro-2-hydroxy-4*H*-1,4,2-oxazaphosphorin-4-yl)-methyl)phosphonate, *N*-oxide, *P*-oxide	417-540-1	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
015-189-00-5	phenyl bis(2,4,6-trimethylbenzoyl)-phosphine oxide	423-340-5	162881-26-7	R43 R53	Xi R: 43-53 S: (2-)22-24-37-61		
016-001-00-4	hydrogen sulphide	231-977-3	7783-06-4	F+; R12 T+; R26 N; R50	F+; T+; N R: 12-26-50 S: (1/2-)9-16-36-38-45-61		

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016-002-00-X	barium sulphide	244-214-4	21109-95-5	R31 Xn; R20/22 N; R50	Xn; N R: 20/22-31-50 S: (2-)28-61		
016-003-00-5	barium polysulphides	256-814-3	50864-67-0	R31 Xi; R36/37/38 N; R50	Xi; N R: 31-36/37/38-50 S: (2-)28-61		
016-004-00-0	calcium sulphide	243-873-5	20548-54-3	R31 Xi; R36/37/38 N; R50	Xi; N R: 31-36/37/38-50 S: (2-)28-61		
016-005-00-6	calcium polysulphides	215-709-2	1344-81-6	R31 Xi; R36/37/38 N; R50	Xi; N R: 31-36/37/38-50 S: (2-)28-61		
016-006-00-1	dipotassium sulphide; potassium sulphide	215-197-0	1312-73-8	R31 C; R34 N; R50	C; N R: 31-34-50 S: (1/2-)26-45-61		
016-007-00-7	potassium polysulphides	253-390-1	37199-66-9	R31 C; R34 N; R50	C; N R: 31-34-50 S: (1/2-)26-45-61		
016-008-00-2	ammonium polysulphides	232-989-1	9080-17-5	R31 C; R34 N; R50	C; N R: 31-34-50 S: (1/2-)26-45-61	C; R34: C ≥ 5 % Xi; R36/38: 1 % ≤ C < 5 % R31: C ≥ 1 %	
016-009-00-8	disodium sulphide; sodium sulphide	215-211-5	1313-82-2	R31 C; R34 N; R50	C; N R: 31-34-50 S: (1/2-)26-45-61		
016-010-00-3	sodium polysulphides	215-686-9	1344-08-7	T; R25 R31 C; R34 N; R50	T; N R: 25-31-34-50 S: (1/2-)26-36/37/39-45-61		
016-011-00-9	sulphur dioxide	231-195-2	7446-09-5	T; R23 C; R34	T R: 23-34 S: (1/2-)9-26-36/37/39-45	T; R23: C ≥ 20 % Xn; R20: 5 % ≤ C < 20 %	5

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016-012-00-4	disulphur dichloride; sulfur monochloride	233-036-2	10025-67-9	R14 T; R25 Xn; R20 R29 C; R35 N; R50	T; C; N R: 14-20-25-29-35-50 S: (1/2-)26-36/37/39-45-61	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
016-013-00-X	sulphur dichloride	234-129-0	10545-99-0	R14 C; R34 Xi; R37 N; R50	C; N R: 14-34-37-50 S: (1/2-)26-45-61		
016-014-00-5	sulphur tetrachloride	—	13451-08-6	R14 C; R34 N; R50	C; N R: 14-34-50 S: (1/2-)26-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
016-015-00-0	thionyl dichloride; thionyl chloride	231-748-8	7719-09-7	R14 Xn; R20/22 R29 C; R35	C R: 14-20/22-29-35 S: (1/2-)26-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
016-016-00-6	sulphuryl chloride	232-245-6	7791-25-5	R14 C; R34 Xi; R37	C R: 14-34-37 S: (1/2-)26-45		
016-017-00-1	chlorosulphonic acid	232-234-6	7790-94-5	R14 C; R35 Xi; R37	C R: 14-35-37 S: (1/2-)26-45		
016-018-00-7	fluorosulphonic acid	232-149-4	7789-21-1	Xn; R20 C; R35	C R: 20-35 S: (1/2-)26-45		
016-019-00-2	oleum ... % SO_3	—	—	R14 C; R35 Xi; R37	C R: 14-35-37 S: (1/2-)26-30-45		B
016-020-00-8	sulphuric acid ... %	231-639-5	7664-93-9	C; R35	C R: 35 S: (1/2-)26-30-45	C; R35: C ≥ 15 % Xi; R36/38: 5 % ≤ C < 15 %	B

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016-021-00-3	methanethiol; methyl mercaptan	200-822-1	74-93-1	F+; R12 T; R23 N; R50-53	F+; T; N R: 12-23-50/53 S: (2-)16-25-60-61		
016-022-00-9	ethanethiol; ethyl mercaptan	200-837-3	75-08-1	F; R11 Xn; R20 N; R50-53	F; Xn; N R: 11-20-50/53 S: (2-)16-25-60-61		
016-023-00-4	dimethyl sulphate	201-058-1	77-78-1	Carc. Cat. 2; R45 Muta. Cat. 3; R68 T+; R26 T; R25 C; R34 R43	T+ R: 45-25-26-34-43-68 S: 53-45	Car. Cat. 2; R45: C ≥ 0,01 % Muta. Cat. 3, R68: C ≥ 0,01 % C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	E
016-024-00-X	dimexano (ISO); bis(methoxythiocarbonyl) disulphide	215-993-8	1468-37-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
016-025-00-5	disul (ISO); 2-(2,4-dichlorophenoxy)ethyl hydrogensulphate; 2,4-DES	205-259-5	149-26-8	Xn; R22 Xi; R38-41	Xn R: 22-38-41 S: (2-)26		
016-026-00-0	sulphamidic acid; sulphamic acid; sulfamic acid	226-218-8	5329-14-6	Xi; R36/38 R52-53	Xi R: 36/38-52/53 S: (2-)26-28-61		
016-027-00-6	diethyl sulphate	200-589-6	64-67-5	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Xn; R20/21/22 C; R34	T R: 45-46-20/21/22-34 S: 53-45		E
016-028-00-1	sodium dithionite; sodium hydrosulphite	231-890-0	7775-14-6	R7 R31 Xn; R22	Xn R: 7-22-31 S: (2-)7/8-26-28-43		
016-029-00-7	*p*-toluenesulphonic acid, containing more than 5 % H_2SO_4	—	—	C; R34	C R: 34 S: (1/2-)26-37/39-45	C; R34: C ≥ 25 % Xi; R36/38: 10 % ≤ C < 25 %	

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016-030-00-2	*p*-toluenesulphonic acid (containing a maximum of 5 % H_2SO_4)	203-180-0	104-15-4	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)26-37		
016-031-00-8	tetrahydrothiophene-1,1-dioxide; sulpholane	204-783-1	126-33-0	Xn; R22	Xn R: 22 S: (2-)25		
016-032-00-3	1,3-propanesultone; 1,2-oxathiolane 2,2-dioxide	214-317-9	1120-71-4	Carc. Cat. 2; R45 Xn; R21/22	T R: 45-21/22 S: 53-45	Carc. Cat. 2; R45: C ≥ 0,01 %	E
016-033-00-9	dimethylsulfamoylchloride	236-412-4	13360-57-1	Carc. Cat. 2; R45 T+; R26 Xn; R21/22 C; R34	T+ R: 45-21/22-26-34 S: 53-45		E
016-034-00-4	tetrasodium 3,3'-(piperazine-1,4-diylbis((6-chloro-1,3,5-triazine-2,4-diyl)imino(2-acetamido)-4,1-phenyleneazo))bis (naphthalene-1,5-disulphonate)	400-010-9	81898-60-4	R43	Xi R: 43 S: (2-)22-24-37		
016-035-00-X	pentasodium 5-anilino-3-(4-(4-(6-chloro-4-(3-sulphonatoanilino)-1,3,5-triazin-2-ylamino)-2,5-dimethylphenylazo)-2,5-disulphonatophenylazo)-4-hydroxynaphthalene-2,7-disulphonate	400-120-7	—	Xi; R36	Xi R: 36 S: (2-)22-26		
016-036-00-5	tetrasodium 5-(4,6-dichloro-5-cyanopyrimidin-2-ylamino)-4-hydroxy-2,3-azodinaphthalene-1,2,5,7-disulphonate	400-130-1	—	R42 N; R51-53	Xn; N R: 42-51/53 S: (2-)22-61		
016-037-00-0	disodium 1-amino-4-(4-benzenesulphonamido-3-sulphonatoanilino)anthraquinone-2-sulphonate	400-350-8	85153-93-1	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
016-038-00-6	disodium 6-((4-chloro-6-(*N*-methyl)-2-toluidino)-1,3,5-triazin-2-ylamino)-1-hydroxy-2-(4-methoxy-2-sulphonatophenylazo)naphthalene-3-sulphonate	400-380-1	86393-35-3	R43	Xi R: 43 S: (2-)22-24-37		
016-039-00-1	tetrasodium 2-(6-chloro-4-(4-(2,5-dimethyl-4-(2,5-disulphonatophenylazo)phenylazo)-3-ureidoanilino)-1,3,5-triazin-2-ylamino)benzene-1,4-disulphonate	400-430-2	—	R43	Xi R: 43 S: (2-)22-24-37		

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016-040-00-7	reaction mass of disodium 6-(2,4-dihydroxyphenylazo)-3-(4-(4-(2,4-dihydroxyphenylazo)anilino)-3-sulphonatophenylazo)-4-hydroxynaphthalene-2-sulphonate and disodium 6-(2,4-diaminophenylazo)-3-(4-(4-(2,4-diaminophenylazo)anilino)-3-sulphonatophenylazo)-4-hydroxynaphthalene-2-sulphonate and trisodium 6-(2,4-dihydroxyphenylazo)-3-(4-(4-(7-(2,4-dihydroxyphenylazo)-1-hydroxy-3-sulphonato-2-naphthylazo)anilino)-3-sulphonatophenylazo)-4-hydroxynaphthalene-2-sulphonate	400-570-4	—	Xi; R36	Xi R: 36 S: (2-)26		
016-041-00-2	calcium 2,5-dichloro-4-(4-((5-chloro-4-methyl-2-sulphonatophenyl)azo)-5-hydroxy-3-methylpyrazol-1-yl)benzenesulphonate	400-710-4	—	Xn; R20	Xn R: 20 S: (2-)		
016-042-00-8	tetrasodium 5-benzamido-3-(5-(4-fluoro-6-(1-sulphonato-2-naphthylamino)-1,3,5-triazin-2-ylamino)-2-sulphonatophenylazo)-4-hydroxynaphthalene-2,7- disulphonate	400-790-0	85665-97-0	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)22-24/25-37		
016-043-00-3	dilithium 6-acetamido-4-hydroxy-3-(4-((2-sulphonatooxy)ethylsulphonyl)phenylazo)naphthalene-2-sulphonate	401-010-1	—	R43	Xi R: 43 S: (2-)24-37		
016-044-00-9	disodium *S,S*-hexane-1,6-diyldi(thiosulphate) dihydrate	401-320-7	—	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
016-045-00-4	lithium sodium hydrogen 4-amino-6-(5-(5-chloro-2,6-difluoropyrimidin-4-ylamino)-2-sulphonatophenylazo)-5-hydroxy-3-(4-(2-(sulphonatooxy)ethylsulphonyl)phenylazo)naphthalene-2,7-disulphonate	401-560-2	108624-00-6	R43	Xi R: 43 S: (2-)22-24-37		
016-046-00-X	sodium hydrogensulphate	231-665-7	7681-38-1	Xi; R41	Xi R: 41 S: (2-)24-26		
016-047-00-5	hexasodium 7-(4-(4-(4-(2,5-disulphonatoanilino)-6-fluoro-1,3,5-triazin-2-ylamino)-2-methylphenylazo)-7-sulphonato-naphthylazo)naphthalene-1,3,5- trisulphonate	401-650-1	85665-96-9	R43	Xi R: 43 S: (2-)22-24-37		
016-048-00-0	sodium 3,5-dichloro-2-(5-cyano-2,6-bis(3-hydroxypropylamino)-4-methylpyridin-3-ylazo)benzenesulphonate	401-870-8	—	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-61		

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016-049-00-6	calcium octadecylxylenesulphonate	402-040-8	—	C; R34 N; R51-53	C; N R: 34-51/53 S: (1/2-)26-28-36/37/39-45-61		
016-050-00-1	potassium sodium 5-(4-chloro-6-(*N*-(4-(4-chloro-6-(5-hydroxy-2,7-disulphonato-6-(2-sulphonatophenylazo)-4-naphthylamino)-1,3,5-triazin-2-ylamino) phenyl-N-methyl) amino)-1,3,5-triazin-2-ylamino)-4-hydroxy-3-(2-sulphonato-phenylazo)naphthalene-2,7-disulphonat	402-150-6	—	Xi; R36 R43	Xi R: 36-43 S: (2-)22-24-26-37		
016-051-00-7	trisodium 7-(4-(6-fluoro-4-(2-(2-vinylsulphonylethoxy)ethylamino)-1,3,5-triazin-2-ylamino)-2-ureidophenylazo)naphthalene-1,3,6- trisulphonate	402-170-5	106359-91-5	R43	Xi R: 43 S: (2-)22-24-37		
016-052-00-2	benzyltributylammonium 4-hydroxynaphthalene-1-sulphonate	402-240-5	102561-46-6	Xn; R20 N; R51-53	Xn; N R: 20-51/53 S: (2-)22-61		
016-053-00-8	(C_{16} or C_{18}-n-alkyl)(C_{16} or C_{18}-n-alkyl)ammonium 2-((C_{16} or C_{18}-n-alkyl)(C_{16} or C_{18}-n-alkyl)carbamoyl)benzenesulphonate	402-460-1	—	Xi; R38 R43 R53	Xi R: 38-43-53 S: (2-)24-37-61		
016-054-00-3	sodium 4-(2,4,4-trimethylpentylcarbonyloxy)benzenesulfonate	400-030-8	—	T; R23-48/23 Xn; R22 Xi; R36/37 R43	T R: 22-23-36/37-43-48/23 S: (1/2-)22-24-36-45		
016-055-00-9	tetrasodium 4-amino-3,6-bis(5-(6-chloro-4-(2-hydroxyethylamino)-1,3,5-triazin-2-ylamino)-2-sulfonatophenylazo)-5-hydroxynaphthalene-2,7-sulfonate (containing > 35 % sodium chloride and sodium acetate)	400-510-7	—	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
016-056-00-4	potassium hydrogensulphate	231-594-1	7646-93-7	C; R34 Xi; R37	C R: 34-37 S: (1/2-)26-36/37/39-45		
016-057-00-X	styrene-4-sulfonyl chloride	404-770-2	2633-67-2	Xi; R38-41 R43	Xi R: 38-41-43 S: (2-)24-26-37/39		
016-058-00-5	thionyl chloride, reaction products with 1,3,4-thiadiazol-2,5-dithiol, *tert*-nonanethiol and C_{12-14}-*tert*-alkylamine	404-820-3	—	Xi; R38 R43 R52-53	Xi R: 38-43-52/53 S: (2-)36/37-61		

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016-059-00-0	*N*,*N*,*N*',*N*'-tetramethyldithiobis(ethylene)diamine dihydrochloride	405-300-9	17339-60-5	Xn; R22 Xi; R36 R43 N; R50-53	Xn; N R: 22-36-43-50/53 S: (2-)26-36/37-60-61		
016-060-00-6	diammonium peroxodisulphate; ammonium persulphate	231-786-5	7727-54-0	O; R8 Xn; R22 Xi; R36/37/38 R42/43	O; Xn R: 8-22-36/37/38-42/43 S: (2-)22-24-26-37		
016-061-00-1	dipotassium peroxodisulphate; potassium persulphate	231-781-8	7727-21-1	O; R8 Xn; R22 Xi; R36/37/38 R42/43	O; Xn R: 8-22-36/37/38-42/43 S: (2-)22-24-26-37		
016-062-00-7	bensultap (ISO); 1,3-bis(phenylsulfonylthio)-2-(*N*,*N*-dimethylamino)propane	—	17606-31-4	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
016-063-00-2	sodium metabisulphite	231-673-0	7681-57-4	Xn; R22 Xi; R41 R31	Xn R: 22-31-41 S: (2-)26-39-46		
016-064-00-8	sodium hydrogensulphite … %; sodium bisulphite … %	231-548-0	7631-90-5	Xn; R22 R31	Xn R: 22-31 S: (2-)25-46		B
016-065-00-3	sodium 1-amino-4-[2-methyl-5-(4-methylphenylsulfonylamino)phenylamino]anthraquinone-2-sulfonate	400-100-8	84057-97-6	N; R51-53	N R: 51/53 S: 61		
016-066-00-9	tetrasodium [5-((4-amino-6-chloro-1,3,5-triazin-2-yl)amino)-2-((2-hydroxy-3,5-disulfonatophenylazo)-2- sulfonatobenzylidenehydrazino)benzoate]copper(II)	404-070-7	116912-62-0	R52-53	R: 52/53 S: 61		
016-067-00-4	(4-methylphenyl)mesitylene sulfonate	407-530-5	67811-06-7	R53	R: 53 S: 61		
016-068-00-X	sodium 3,5-bis(tetradecyloxycarbonyl)benzenesulfinate	407-720-8	155160-86-4	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		

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016-069-00-5	3,5-bis-(tetradecyloxycarbonyl)benzenesulfinic acid	407-990-7	141915-64-2	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
016-070-00-0	4-benzyloxy-4'-(2,3-epoxy-2-methylprop-1-yloxy)diphenyl-sulfone	408-220-2	—	R53	R: 53 S: 61		
016-071-00-6	trisodium 3-amino-6,13-dichloro-10-((3-((4-chloro-6-(2-sulfophenylamino)-1,3,5-triazin-2-yl)amino)propyl) amino)-4,11-triphenoxydioxazinedisulfonate	410-130-3	136248-03-8	R43	Xi R: - 43 S: (2-)22-24-37		
016-072-00-1	3-amino-4-hydroxy-*N*-(2-methoxyethyl)-benzenesulfonamide	411-520-6	112195-27-4	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24-26-37/39-61		
016-073-00-7	tetrakis(phenylmethyl)thioperoxydi(carbothioamide)	404-310-0	10591-85-2	R53	R: 53 S: 61		
016-074-00-2	6-fluoro-2-methyl-3-(4-methylthiobenzyl)indene	405-410-7	—	Xi; R38-41 R43 N; R51-53	Xi; N R: 38-41-43-51/53 S: (2-)26-36/37/39-61		
016-075-00-8	2,2'-diallyl-4,4'-sulfonyldiphenol	411-570-9	41481-66-7	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
016-076-00-3	2,3-bis((2-mercaptoethyl)thio)-1-propanethiol	411-290-7	131538-00-6	Xn; R22-48/22 N; R50-53	Xn; N R: 22-48/22-50/53 S: (2-)23-24/25-36-60-61		
016-077-00-9	2-chloro-*p*-toluenesulfochloride	412-890-1	42413-03-6	C; R34 R43 R52-53	C R: 34-43-52/53 S: (1/2-)23-26-36/37/39-45-61		
016-078-00-4	4-methyl-*N*,*N*-bis(2-(((4-methylphenyl)sulfonyl)amino)ethyl) benzenesulfonamide	413-300-5	56187-04-3	R53	R: 53 S: 61		
016-079-00-X	*N*,*N*-bis(2-(*p*-toluenesulfonyloxy)ethyl)-*p*-toluenesulfonamide	412-920-3	16695-22-0	R43 R53	Xi R: 43-53 S: (2-)24-37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
016-080-00-5	sodium 2-anilino-5-(2-nitro-4-(*N*-phenylsulfamoyl))anilino-benzenesulfonate	412-320-1	31361-99-6	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
016-081-00-0	hexahydrocyclopenta[*c*]pyrrole-1-(1*H*)-ammonium *N*-ethoxy-carbonyl-*N*-(*p*-tolylsulfonyl)azanide	418-350-1	—	Muta. Cat. 3; R68 Xn; R22 Xi; R36 R43 N; R51-53	Xn; N R: 22-36-43-68-51/53 S: (2-)26-36/37-61		
016-082-00-6	ethoxysulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-(2-ethoxyphenoxysulfonyl)urea	—	126801-58-9	N; R50-53	N R: 50/53 S: 60-61		
016-083-00-1	acibenzolar-*S*-methyl; benzo[1,2,3]thiadiazole-7-carbothioic acid *S*-methyl ester	420-050-0	135158-54-2	Xi; R36/37/38 R43 N; R50-53	Xi; N R: 36/37/38-43-50/53 S: (2-)24/25-37-46-59-60-61		
016-084-00-7	prosulfuron; 1-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)-3-[2-(3,3,3-trifluoropropyl)phenylsulfonyl]urea	—	94125-34-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
016-085-00-2	flazasulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-(3-trifluoromethyl-2-pyridylsulfonyl)urea	—	104040-78-0	N; R50-53	N R: 50/53 S: 60-61		
016-086-00-8	tetrasodium 10-amino-6,13-dichloro-3-(3-(4-(2,5-disulfonatoanilino)-6-fluoro-1,3,5-triazin-2-ylamino)prop-3-ylamino)-5,12-dioxa-7,14-diazapentacene-4,11-disulfonate	402-590-9	109125-56-6	Xi; R41	Xi R: 41 S: (2-)22-26-39		

016-087-00-3	reaction mass of: thiobis(4,1-phenylene)-*S*,*S*,*S*',*S*'-tetraphenyldisulfonium bishexafluorophosphate;
	diphenyl(4-phenylthiophenyl)sulfonium hexafluorophosphate;
	propylene carbonate
		403-490-8	104558-95-4	Xi; R36
				R43
				N; R50-53
					Xi; N
					R: 36-43-50/53
					S: (2-)24-26-37-60-61		

016-088-00-9	4-(bis(4-(diethylamino)phenyl)methyl)benzene-1,2-dimethanesulfonic acid	407-280-7	71297-11-5	R52-53	R: 52/53 S: 61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
016-089-00-4	reaction mass of esters of 5,5',6,6',7,7'-hexahydroxy-3,3,3',3'-tetramethyl-1,1'-spirobiindan and 2-diazo-1,2-dihydro-1-oxo-5-sulfonaphthalene	413-840-1	—	E; R2 F; R11 R53	E R: 2-11-53 S: (2-)33-35-40-61		
016-090-00-X	4-methyl-*N*-(methylsulfonyl)benzenesulfonamide	415-040-8	14653-91-9	Xn; R22 Xi; R37-41	Xn R: 22-37-41 S: (2-)26-39		
016-091-00-5	C_{12-14}-*tert*-alkyl ammonium 1-amino-9,10-dihydro-9,10-dioxo-4-(2,4,6-trimethylanilino)-anthracen-2-sulfonate	414-110-5	—	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
016-093-00-6	reaction mass of: 4-(7-hydroxy-2,4,4-trimethyl-2-chromanyl) resorcinol-4-yl-tris(6-diazo-5,6-dihydro-5-oxonaphthalen-1-sulfonate); 4-(7-hydroxy-2,4,4-trimethyl-2-chromanyl)resorcinolbis(6-diazo-5,6-dihydro-5-oxonaphthalen-1-sulfonate) (2:1)	414-770-4	140698-96-0	F; R11 Carc. Cat. 3; R40	F; Xn R: 11-40 S: (2-)7-36/37		
016-095-00-7	reaction mass of: reaction product of 4,4'-methylenebis[2-(4-hydroxybenzyl)-3,6-dimethylphenol] and 6-diazo-5,6-dihydro-5-oxo-naphthalenesulfonate (1:2); Reaction product of 4,4'-methylenebis[2-(4-hydroxybenzyl)-3,6-dimethylphenol] and 6-diazo-5,6-dihydro-5-oxo-naphthalenesulfonate (1:3)	417-980-4	—	F; R11 Carc. Cat. 3; R40	F; Xn R: 11-40 S: (2-)7-36/37		
016-096-00-2	thifensulfuron-methyl (ISO); methyl 3-(4-methoxy-6-methyl-1,3,5-triazin-2-ylcarbamoylsulfamoyl)thiophene-2-carboxylate	—	79277-27-3	N; R50-53	N R: 50/53 S: 60-61		
017-001-00-7	chlorine	231-959-5	7782-50-5	T; R23 Xi; R36/37/38 N; R50	T; N R: 23-36/37/38-50 S: (1/2-)9-45-61		
017-002-00-2	hydrogen chloride	231-595-7	7647-01-0	T; R23 C; R35	T; C R: 23-35 S: (1/2-)9-26-36/37/39-45		5
017-002-01-X	hydrochloric acid ... %	231-595-7	—	C; R34 Xi; R37	C R: 34-37 S: (1/2-)26-45	C; R34-37: C ≥ 25 % Xi; R36/37/38: 10 % ≤ C < 25 %	B

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
017-003-00-8	barium chlorate	236-760-7	13477-00-4	O; R9 Xn; R20/22 N; R51-53	O; Xn; N R: 9-20/22-51/53 S: (2-)13-27-61		
017-004-00-3	potassium chlorate	223-289-7	3811-04-9	O; R9 Xn; R20/22 N; R51-53	O; Xn; N R: 9-20/22-51/53 S: (2-)13-16-27-61		
017-005-00-9	sodium chlorate	231-887-4	7775-09-9	O; R9 Xn; R22 N; R51-53	O; Xn; N R: 9-22-51/53 S: (2-)13-17-46-61		
017-006-00-4	perchloric acid ... %	231-512-4	7601-90-3	R5 O; R8 C; R35	O; C R: 5-8-35 S: (1/2-)23-26-36-45	C; R35: C ≥ 50 % C; R34: 10 % ≤ C < 50 % Xi; R36/38: 1 % ≤ C < 10 % Footnote O; R5-8: C > 50 %	B
017-007-00-X	barium perchlorate	236-710-4	13465-95-7	O; R9 Xn; R20/22	O; Xn R: 9-20/22 S: (2-)27		
017-008-00-5	potassium perchlorate	231-912-9	7778-74-7	O; R9 Xn; R22	O; Xn R: 9-22 S: (2-)13-22-27		
017-009-00-0	ammonium perchlorate	232-235-1	7790-98-9	O; R9 R44 ⊗	O R: 9-44 S: (2-)14-16-27-36/37		G
017-010-00-6	sodium perchlorate	231-511-9	7601-89-0	O; R9 Xn; R22	O; Xn R: 9-22 S: (2-)13-22-27		
017-011-00-1	sodium hypochlorite, solution ... % Cl active	231-668-3	7681-52-9	C; R34 R31 N; R50	C; N R: 31-34-50 S: (1/2-)28-45-50-61	R31: C ≥ 5 %	B
017-012-00-7	calcium hypochlorite	231-908-7	7778-54-3	O; R8 ⊗ Xn; R22 R31 C; R34 N; R50	O; C; N R: 8-22-31-34-50 S: (1/2-)26-36/37/39-45-61	C; R34: C ≥ 10 % Xi; R37/38-41: 3 % ≤ C < 10 % Xi; R36: 0,5 % ≤ C < 3 %	

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017-013-00-2	calcium chloride	233-140-8	10043-52-4	Xi; R36	Xi R: 36 S: (2-)22-24		
017-014-00-8	ammonium chloride	235-186-4	12125-02-9	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)22		
017-015-00-3	(2-(aminomethyl)phenyl)acetylchloride hydrochloride	417-410-4	61807-67-8	Xn; R22 C; R35 R43	C R: 22-35-43 S: (1/2-)26-36/37/39-45		
017-016-00-9	methyltriphenylphosphonium chloride	418-400-2	1031-15-8	Xn; R21/22 Xi; R38-41 N; R51-53	Xn; N R: 21/22-38-41-51/53 S: (2-)22-26-36/37/39-61		
017-017-00-4	(Z)-13-docosenyl-*N*,*N*-bis(2-hydroxyethyl)-*N*-methyl-ammonium-chloride	426-210-6	120086-58-0	C; R34 N; R50-53	C; N R: 34-50/53 S: (2-)26-36/37/39-45-60-61		
017-018-00-X	*N*,*N*,*N*-trimethyl-2,3-bis(stearoyloxy)propylammonium chloride	405-660-7	—	N; R51-53	N R: 51/53 S: 61		
017-019-00-5	(*R*)-1,2,3,4-tetrahydro-6,7-dimethoxy-1-veratrylisoquinoline hydrochloride	415-110-8	54417-53-7	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
017-020-00-0	ethyl propoxy aluminium chloride	421-790-7	13014-29-4	F; R15 R 14 C; R35	F; C R: 14/15-35 S: (1/2-)16-23-26-30-36/37/39-43-45		
017-021-00-6	behenamidopropyl-dimethyl-(dihydroxypropyl) ammonium chloride	423-420-1	136920-10-0	Xi; R41 R43 N; R50-53	Xi; N R: 41-43-50/53 S: (2-)26-36/37/39-60-61		
019-001-00-2	potassium	231-119-8	7440-09-7	F; R15 R14 C; R34	F; C R: 14/15-34 S: (1/2-)5-8-45		

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019-002-00-8	potassium hydroxide; caustic potash	215-181-3	1310-58-3	Xn; R22 C; R35	C R: 22-35 S: (1/2-)26-36/37/39-45	C; R35: C ≥ 5 % C; R34: 2 % ≤ C < 5 % Xi; R36/38: 0.5 % ≤ C < 2 %	
020-001-00-X	calcium	231-179-5	7440-70-2	F; R15	F R: 15 S: (2-)8-24/25-43		
020-002-00-5	calcium cyanide	209-740-0	592-01-8	T+; R28 R32 N; R50-53	T+; N R: 28-32-50/53 S: (1/2-)7/8-23-36/37-45-60-61		

020-003-00-0	reaction mass of: dicalcium (bis(2-hydroxy-5-tetra-propenyl-phenylmethyl)methylamine)dihydroxide;
	tri-calcium (tris(2-hydroxy-5-tetra-propenylphenylmethyl) methylamine)tri-hydroxide;
	poly[calcium ((2-hydroxy-5-tetra-propenyl-phenylmethyl) methylamine)hydroxide]
		420-470-4	—	Xi; R36/38
				R43
					Xi
					R: 36/38-43
					S: (2-)24-26-37		

022-001-00-5	titanium tetrachloride	231-441-9	7550-45-0	R14 C; R34	C R: 14-34 S: (1/2-)7/8-26-36/37/39-45		
022-002-00-0	titanium(4+) oxalate	403-260-7	—	Xi; R41	Xi R: 41 S: (2-)26-39		
022-003-00-6	bis(η^5-cyclopentadienyl)-bis(2,6-difluoro-3-[pyrrol-1-yl]-phenyl)titanium	412-000-1	125051-32-3	F; R11 Repr. Cat. 3; R62 Xn; R48/22 N; R51-53	F; Xn; N R: 11-48/22-62-51/53 S: (2-)7-22-33-36/37-61		
023-001-00-8	divanadium pentaoxide; vanadium pentoxide	215-239-8	1314-62-1	Muta. Cat. 3; R68 Repr. Cat. 3; R63 T; R48/23 Xn; R20/22 Xi; R37 N; R51-53	T; N R: 20/22-37-48/23-51/53-63-68 S: (1/2-)36/37-38-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
024-001-00-0	chromium (VI) trioxide	215-607-8	1333-82-0	O; R9 Carc. Cat. 1; R45 Muta. Cat. 2; R46 Repr. Cat. 3; R62 T+; R26 T; R24/25-48/23 C; R35 R42/43 N; R50-53	O; T+; N R: 45-46-9-24/25-26-35-42/43-48/23-62-50/53 S: 53-45-60-61	C: R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	E
024-002-00-6	potassium dichromate	231-906-6	7778-50-9	O; R8 Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23 Xn; R21 C; R34 R42/43 N; R50-53	T+; N; O R: 45-46-60-61-8-21-25-26-34-42/43-48/23-50/53 S: 53-45-60-61	C; R34: C ≥ 10 %: Xi; R36/37/38: 5 % ≤ C < 10 %	E 3
024-003-00-1	ammonium dichromate	232-143-1	7789-09-5	E; R2 O; R8 Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23 Xn; R21 C; R34 R42/43 N; R50-53	E; T+; N R: 45-46-60-61-2-8-21-25-26-34-42/43-48/23-50/53 S: 53-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 % R42/43: C ≥ 0,2 %	E 3

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024-004-00-7	sodium dichromate anhydrate	234-190-3	10588-01-9	O; R8 Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23 Xn; R21 C; R34 R42/43 N; 50-53	T+; N; O R: 45-46-60-61-8-21-25-26-34-42/43-48/23-50/53 S: 53-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 % R42/43: C ≥ 0,2 %	E 3
024-004-01-4	sodium dichromate, dihydrate	234-190-3	7789-12-0	O; R8 Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23 Xn; R21 C; R34 R42/43 N; R50-53	T+; N; O R: 45-46-60-61-8-21-25-26-34-42/43-48/23-50/53 S: 53-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 % R42/43: C ≥ 0,2 %	E 3
024-005-00-2	chromyl dichloride; chromic oxychloride	239-056-8	14977-61-8	O; R8 Carc. Cat. 2; R49 Muta. Cat. 2; R46 C; R35 R43 N; R50-53	O; T; C; N R: 49-46-8-35-43-50/53 S: 53-45-60-61	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 0,5 % ≤ C < 5 % R43: C ≥ 0,5 %	3
024-006-00-8	potassium chromate	232-140-5	7789-00-6	Carc. Cat. 2; R49 Muta. Cat. 2; R46 Xi; R36/37/38 R43 N; R50-53	T; N R: 49-46-36/37/38-43-50/53 S: 53-45-60-61	R43: C ≥ 0,5 %	3

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024-007-00-3	zinc chromates including zinc potassium chromate	—	—	Carc. Cat. 1; R45 Xn; R22 R43 N; R50-53	T; N R: 45-22-43-50/53 S: 53-45-60-61		AE
024-008-00-9	calcium chromate	237-366-8	13765-19-0	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		E
024-009-00-4	strontium chromate	232-142-6	7789-06-2	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		E
024-010-00-X	dichromium tris(chromate); chromium III chromate; chromic chromate	246-356-2	24613-89-6	O; R8 Carc. Cat. 2; R45 C; R35 R43 N; R50-53	O; T; C; N R: 45-8-35-43-50/53 S: 53-45-60-61		
024-011-00-5	ammonium bis(1-(3,5-dinitro-2-oxidophenylazo)-3-(*N*-phenylcarbamoyl)-2-naphtholato)chromate(1-)	400-110-2	109125-51-1	F; R11 N; R50-53	F; N R: 11-50/53 S: (2-)33-60-61		
024-012-00-0	trisodium bis(7-acetamido-2-(4-nitro-2-oxidophenylazo)-3-sulphonato-1-naphtholato)chromate(1-)	400-810-8	—	Muta. Cat. 3; R68	Xn R: 68 S: (2-)22-36/37		
024-013-00-6	trisodium (6-anilino-2-(5-nitro-2-oxidophenylazo)-3-sulphonato-1-naphtholato)(4-sulphonato-1,1'-azodi-2,2'naphtholato)chromate(1-)	402-500-8	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
024-014-00-1	trisodium bis(2-(5-chloro-4-nitro-2-oxidophenylazo)-5-sulphonato-1-naphtholato)chromate(1-)	402-870-0	93952-24-0	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
024-015-00-7	disodium (3-methyl-4-(5-nitro-2-oxidophenylazo)-1-phenylpyrazololato)(1-(3-nitro-2-oxido-5-sulfonatophenylazo)-2-naphtholato)chromate(1-)	404-930-1	—	Xn; R20 Xi; R41 N; R51-53	Xn; N R: 20-41-51/53 S: (2-)26-39-61		

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024-016-00-2	tetradecylammonium bis(1-(5-chloro-2-oxidophenylazo)-2-naphtholato)chromate(1-)	405-110-6	88377-66-6	Xn; R48/22 R53	Xn R: 48/22-53 S: (2-)22-36-61		
024-017-00-8	Chromium (VI) compounds, with the exception of barium chromate and of compounds specified elsewhere in this Annex	—	—	Carc. Cat. 2; R49 R43 N; R50-53	T; N R: 49-43-50/53 S: 53-45-60-61		A
024-018-00-3	sodium chromate	231-889-5	7775-11-3	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23 Xn; R21 C; R34 R42/43 N; R50-53	T+; N R: 45-46-60-61-21-25-26-34-42/43-48/23-50/53 S: 53-45-60-61	R42/43: C ≥ 0,2 %	E 3
024-019-00-9	Main component: acetoacetic acid anilide/3-amino-1-hydroxybenzene (ATAN-MAP): trisodium {6-[(2 or 3 or 4)-amino-(4 or 5 or 6)-hydroxyphenylazo]-5'-(phenylsulfamoyl)-3-sulfonatonaphthalene-2-azobenzene-1,2'-diolato}-{6"-[1-(phenylcarbamoyl)ethylazo]-5'"-(phenylsulfamoyl)-3"-sulfonatonaphthalene-2"-azobenzene-1",2'"-diolato}chromate (III); by-product 1: acetoacetic acid anilide/acetoacetic acid anilide (ATAN-ATAN): trisodium bis{6-[1-(phenylcarbamoyl)ethylazo]-5'-(phenylsulfonyl)-3-sulfonatonaphthalene-2-azobenzene-1,2'-diolato}chromate (III); by-product 2: 3-amino-1-hydroxybenzene/3-amino-1-hydroxybenzene (MAP-MAP): trisodium bis{6-[(2 or 3 or 4)-amino-(4 or 5 or 6)-hydroxyphenylazo]-5'-(phenylsulfamoyl)-3-sulfonatonaphthalene-2-azobenzene-1,2'-diolato} chromate (III)	419-230-1	—	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
024-020-00-4	trisodium bis[(3'-nitro-5'-sulfonato(6-amino-2-[4-(2-hydroxy-1-naphtylazo)phenylsulfonylamino]pyrimidin-5-azo)benzene-2',4-diolato)]chromate (III)	418-220-4	—	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
025-001-00-3	manganese dioxide	215-202-6	1313-13-9	Xn; R20/22	Xn R: 20/22 S: (2-)25		

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025-002-00-9	potassium permanganate	231-760-3	7722-64-7	O; R8 Xn; R22 N; R50-53	O; Xn; N R: 8-22-50/53 S: (2-)60-61		
025-003-00-4	manganese sulphate	232-089-9	7785-87-7	Xn; R48/20/22 N; R51-53	Xn; N R: 48/20/22-51/53 S: (2-)22-61		
025-004-00-X	bis(*N*,*N*',*N*''-trimethyl-1,4,7-triazacyclononane)-trioxo-dimanganese (IV) di(hexafluorophosphate) monohydrate	411-760-1	116633-53-5	N; R51-53	N R: 51/53 S: 61		
025-005-00-5	reaction mass of: tri-sodium [29*H*, 31*H*-phthalocyanine-*C*,*C*,*C*-trisulfonato (6-)-*N*29,*N*30,*N*31,*N*32] manganate (3-); tetrasodium [29*H*,31*H*-phthalocyanine-*C*,*C*,*C*,*C*-tetrasulfonato (6-)-*N*29,*N*30,*N*31,*N*32], manganate (3-); pentasodium [29*H*,31*H*-phthalocyanine-*C*,*C*,*C*,*C*,*C*-pentasulfonato (6-)-*N*29,*N*30,*N*31,*N*32] manganate (3-)	417-660-4	—	N; R50-53	N R: 50/53 S: 60-61		
026-001-00-6	(η-cumene)-(η-cyclopentadienyl)iron(II) hexafluoroantimonate	407-840-0	100011-37-8	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)22-26-39-61		
026-002-00-1	(η-cumene)-(η-cyclopentadienyl)iron(II) trifluoromethane-sulfonate	407-880-9	117549-13-0	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)26-61		
027-001-00-9	cobalt	231-158-0	7440-48-4	R42/43 R53	Xn R: 42/43-53 S: (2-)22-24-37-61		
027-002-00-4	cobalt oxide	215-154-6	1307-96-6	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
027-003-00-X	cobalt sulphide	215-273-3	1317-42-6	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
027-004-00-5	cobalt dichloride	231-589-4	7646-79-9	Carc. Cat. 2; R49 Xn; R22 R42/43 N; R50-53	T; N R: 49-22-42/43-50/53 S: (2-)22-53-45-60-61	Carc. Cat. 2; R49: C ≥ 0,01 % Xn; R22: C ≥ 2,5 %	E 1

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027-005-00-0	cobalt sulphate	233-334-2	10124-43-3	Carc. Cat. 2; R49 Xn; R22 R42/43 N; R50-53	T; N R: 49-22-42/43-50/53 S: (2-)22-53-45-60-61	Carc. Cat. 2; R49: C ≥ 0,01 %	E 1
028-001-00-1	tetracarbonylnickel; nickel tetracarbonyl	236-669-2	13463-39-3	F; R11 Carc. Cat. 3; R40 Repr. Cat. 2; R61 T+; R26 N; R50-53	F; T+; N R: 61-11-26-40-50/53 S: 53-45-60-61		E
028-002-00-7	nickel	231-111-4	7440-02-0	Carc. Cat. 3; R40 R43	Xn R: 40-43 S: (2-)22-36		
028-003-00-2	nickel monoxide	215-215-7	1313-99-1	Carc. Cat. 1; R49 R43 R53	T R: 49-43-53 S: 53-45-61		
028-004-00-8	nickel dioxide	234-823-3	12035-36-8	Carc. Cat. 1; R49 R43 R53	T R: 49-43-53 S: 53-45-61		
028-005-00-3	dinickel trioxide	215-217-8	1314-06-3	Carc. Cat. 1; R49 R43 R53	T R: 49-43-53 S: 53-45-61		
028-006-00-9	nickel sulphide	240-841-2	16812-54-7	Carc. Cat. 1; R49 R43 N; R50-53	T; N R: 49-43-50/53 S: 53-45-60-61		
028-007-00-4	nickel subsulphide; trinickel disulphide	234-829-6	12035-72-2	Carc. Cat. 1; R49 R43 N; R51-53	T; N R: 49-43-51/53 S: 53-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
028-008-00-X	nickel dihydroxide	235-008-5	12054-48-7	Carc. Cat. 3; R40 Xn; R20/22 R43 N; R50-53	Xn; N R: 20/22-40-43-50/53 S: (2-)22-36-60-61		
028-009-00-5	nickel sulphate	232-104-9	7786-81-4	Carc. Cat. 3; R40 Xn; R22 R42/43 N; R50-53	Xn; N R: 22-40-42/43-50/53 S: (2-)22-36/37-60-61		
028-010-00-0	nickel carbonate	222-068-2	3333-67-3	Carc. Cat. 3; R40 Xn; R22 R43 N; R50-53	Xn; N R: 22-40-43-50/53 S: (2-)22-36/37-60-61		
029-001-00-4	copper chloride; copper (I) chloride; cuprous chloride	231-842-9	7758-89-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-60-61		
029-002-00-X	dicopper oxide; copper (I) oxide	215-270-7	1317-39-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-60-61		
029-003-00-5	Naphthenic acids, copper salts; copper naphthenate	215-657-0	1338-02-9	R10 Xn; R22 N; R50-53	Xn; N R: 10-22-50/53 S: (2-)60-61		
029-004-00-0	copper sulphate	231-847-6	7758-98-7	Xn; R22 Xi; R36/38 N; R50-53	Xn; N R: 22-36/38-50/53 S: (2-)22-60-61		
029-005-00-6	(tris(chloromethyl)phthalocyaninato)copper(II), reaction products with *N*-methylpiperazine and methoxyacetic acid	401-260-1	—	Xi; R36	Xi R: 36 S: (2-)26		
029-006-00-1	tris(octadec-9-enylammonium) (trisulfonatophthalocyaninato) copper(II)	403-210-4	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)22-26-39-61		

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029-007-00-7	(trisodium (2-((3-(6-(2-chloro-5-sulfonato)anilino)-4-(3-carboxypyridinio)-1,3,5-triazin-2-ylamino)-2-oxido-5-sulfonatophenylazo)phenylmethylazo)-4-sulfonatobenzoato)copper (3-)) hydroxide	404-670-9	89797-01-3	E; R2 R43	E; Xi R: 2-43 S: (2-)22-24-35-37		
029-008-00-2	copper(II) methanesulfonate	405-400-2	54253-62-2	Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-50/53 S: (2-)26-36/37/39-60-61		
029-009-00-8	phthalocyanine-*N*-[3-(diethylamino)propyl]sulfonamide copper complex	413-650-9	93971-95-0	R52-53	R: 52/53 S: 61		
029-010-00-3	reaction mass of compounds from (dodecakis(*p*-tolylthio) phthalocyaninato)copper(II) to (hexadecakis(*p*-tolylthio) phthalocyaninato)copper(II)	407-700-9	101408-30-4	R43	Xi R: 43 S: (2-)24-37		
029-011-00-9	sodium [29*H*,31*H*-phthalocyaninato-(2-)-*N*29,*N*30,*N*31,*N*32]-((3-(*N*-methyl-*N*-(2-hydroxyethyl)amino)propyl)amino) sulfonyl-sulfonato, copper complex	412-730-0	150522-10-4	C; R34	C R: 34 S: (1/2-)22-26-36/37/39-45		
029-012-00-4	sodium ((*N*-(3-trimethylammoniopropyl)sulfamoyl)methylsulfonatophthalocyaninato)copper(II)	407-340-2	124719-24-0	Xi; R41	Xi R: 41 S: (2-)26-39		
029-013-00-X	trisodium(2-(α-(3-(4-chloro-6-(2-(2-(vinylsulfonyl)ethoxy) ethylamino)-1,3,5-triazin-2-ylamino)-2-oxido-5-sulfonatophenylazo)benzylidenehydrazino)-4-sulfonatobenzoato)copper(II)	407-580-8	130201-51-3	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)24-37-61		
030-001-00-1	zinc powder — zinc dust (pyrophoric)	231-175-3	7440-66-6	F; R15-17 N; R50-53	F; N R: 15-17-50/53 S: (2-)43-46-60-61		
030-001-01-9	zinc powder — zinc dust (stabilised)	231-175-3	7440-66-6	N; R50-53	N R: 50/53 S: 60-61		
030-003-00-2	zinc chloride	231-592-0	7646-85-7	C; R34 Xn; R22 N; R50-53	C; N R: 22-34-50/53 S: (1/2-)26-36/37/39-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	

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030-004-00-8	dimethylzinc; [1] diethylzinc [2]	208-884-1 [1] 209-161-3 [2]	544-97-8 [1] 557-20-0 [2]	F; R17 R14 C; R34 N; R50-53	F; C; N R: 14-17-34-50/53 S: (1/2-)16-43-45-60-61		
030-005-00-3	diamminediisocyanatozinc	401-610-3	—	Xn; R22 Xi; R41 R42/43 N; R50	Xn; N R: 22-41-42/43-50 S: (2-)22-26-36/37/39-41-61		
030-006-00-9	zinc sulphate (hydrous) (mono-, hexa- and hepta hydrate); [1] zinc sulphate (anhydrous) [2]	231-793-3 [1] 231-793-3 [2]	7446-19-7 [1] 7733-02-0 [2]	Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-50/53 S: (2-)22-26-39-46-60-61		
030-007-00-4	bis(3,5-di-*tert*-butylsalicylato-O^1,O^2)zinc	403-360-0	42405-40-3	F; R11 Xn; R22 N; R50-53	F; Xn; N R: 11-22-50/53 S: (2-)7-22-60-61		
030-008-00-X	hydroxo(2-(benzenesulfonamido)benzoato)zinc(II)	403-750-0	113036-91-2	Xn; R20 N; R51-53	Xn; N R: 20-51/53 S: (2-)22-57-61		
030-011-00-6	trizinc bis(orthophosphate)	231-944-3	7779-90-0	N; R50-53	N R: 50/53 S: 60-61		
030-013-00-7	zinc oxide	215-222-5	1314-13-2	N; R50-53	N R: 50/53 S: 60-61		
033-001-00-X	arsenic	231-148-6	7440-38-2	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)20/21-28-45-60-61		
033-002-00-5	arsenic compounds, with the exception of those specified elsewhere in this Annex	—	—	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)20/21-28-45-60-61	T; R23/25: C ≥ 0,2 % Xn; R20/22: 0,1 % ≤ C < 0,2 %	A 1

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033-003-00-0	diarsenic trioxide; arsenic trioxide	215-481-4	1327-53-3	Carc. Cat. 1; R45 T+; R28 C; R34 N; R50-53	T+; N R: 45-28-34-50/53 S: 53-45-60-61		E
033-004-00-6	diarsenic pentaoxide; arsenic pentoxide; arsenic oxide	215-116-9	1303-28-2	Carc. Cat. 1; R45 T; R23/25 N; R50-53	T; N R: 45-23/25-50/53 S: 53-45-60-61		E
033-005-00-1	arsenic acid and its salts	—	—	Carc. Cat. 1; R45 T; R23/25 N; R50-53	T; N R: 45-23/25-50/53 S: 53-45-60-61		AE
033-006-00-7	arsine	232-066-3	7784-42-1	F+; R12 T+; R26 Xn; R48/20 N; R50-53	F+; T+; N R: 12-26-48/20-50/53 S: (1/2-)9-16-28-33-36/37-45-60-61		
033-007-00-2	*tert*-butylarsine	423-320-6	4262-43-5	F; R17 T+; R26	F; T+ R: 17-26 S: (1/2-)9-28-36/37-43-45		
034-001-00-2	selenium	231-957-4	7782-49-2	T; R23/25 R33 R53	T R: 23/25-33-53 S: (1/2-)20/21-28-45-61		
034-002-00-8	selenium compounds except cadmium sulphoselenide	—	—	T; R23/25 R33 N; R50-53	T; N R: 23/25-33-50/53 S: (1/2-)20/21-28-45-60-61		A
034-003-00-3	sodium selenite	233-267-9	10102-18-8	T+; R28 T; R23 R31 R43 N; R51-53	T+; N R: 23-28-31-43-51/53 S: (1/2-)28-36/37-45-61		

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035-001-00-5	bromine	231-778-1	7726-95-6	T+; R26 C; R35 N; R50	T+; C; N R: 26-35-50 S: (1/2-)7/9-26-45-61		
035-002-00-0	hydrogen bromide	233-113-0	10035-10-6	C; R35 Xi; R37	C R: 35-37 S: (1/2-)7/9-26-45		
035-002-01-8	hydrobromic acid ... %	—	—	C; R34 Xi; R37	C R: 34-37 S: (1/2-)7/9-26-45	C; R34: C ≥ 40 % Xi; R36/38: 10 % ≤ C < 40 % Xi; R37: C ≥ 10 %	B
035-003-00-6	potassium bromate	231-829-8	7758-01-2	O; R9 Carc. Cat. 2; R45 T; R25	T; O R: 45-9-25 S: 53-45		E
035-004-00-1	2-hydroxyethylammonium perbromide	407-440-6	—	O; R8 Xn; R22 C; R35 R43 N; R50	O; C; N R: 8-22-35-43-50 S: (1/2-)3/7-14-26-36/37/39-45-60-61		
040-001-00-3	zirconium powder (pyrophoric)	231-176-9	7440-67-7	F; R15-17	F R: 15-17 S: (2-)7/8-43		
040-002-00-9	zirconium powder (non pyrophoric)	—	—	F; R15	F R: 15 S: (2-)7/8-43		
042-001-00-9	molybdenum trioxide	215-204-7	1313-27-5	Xn; R48/20/22 Xi; R36/37	Xn R: 36/37-48/20/22 S: (2-)22-25		
042-002-00-4	tetrakis(dimethylditetradecylammonium) hexa-μ-oxotetra-μ3-oxodi-μ5-oxotetradecaoxooctamolybdate(4-)	404-760-8	117342-25-3	T; R23 Xi; R41 R53	T R: 23-41-53 S: (1/2-)26-37/39-45-61		
042-003-00-X	tetrakis(trimethylhexadecylammonium) hexa-mu-oxotetra-mu3-oxodi-mu5-oxotetradecaoxooctamolybdate(4-)	404-860-1	116810-46-9	F; R11 Xi; R41 N; R50-53	F; Xi; N R: 11-41-50/53 S: (2-)26-39-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
042-004-00-5	Reaction product of ammonium molybdate and C_{12}-C_{24}-diethoxylated alkylamine (1:5-1:3)	412-780-3	—	Xi; R38 R43 N; R51-53	Xi; N R: 38-43-51/53 S: 24/25-37-61		
047-001-00-2	silver nitrate	231-853-9	7761-88-8	⊗ C; R34 N; R50-53	C; N R: 34-50/53 S: (1/2-)26-45-60-61		
048-001-00-5	cadmium compounds, with the exception of cadmium sulphoselenide (xCdS.yCdSe), reaction mass of cadmium sulphide with zinc sulphide (xCdS.yZnS), reaction mass of cadmium sulphide with mercury sulphide (xCdS.yHgS), and those specified elsewhere in this Annex	—	—	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)60-61	Xn; R20/21/22: C ≥ 0,1 %	A 1
048-002-00-0	cadmium (non-pyrophoric); [1] cadmium oxide (non-pyrophoric) [2]	231-152-8 [1] 215-146-2 [2]	7440-43-9 [1] 1306-19-0 [2]	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62-63 T+; R26 T; R48/23/25 N; R50-53	T+; N R: 45-26-48/23/25-62-63-68-50/53 S: 53-45-60-61		E
048-003-00-6	cadmium diformate; cadmiumformate	224-729-0	4464-23-7	T; R23/25 R33 Xn; R68 N; R50-53	T; N R: 23/25-33-68-50/53 S: (1/2-)22-45-60-61	T; R23/25: C ≥ 10 % Xn; R20/22: 0,25 % ≤ C < 10 % R33: C ≥ 0,25 %	
048-004-00-1	cadmium cyanide	208-829-1	542-83-6	T+; R26/27/28 R32 R33 Xn; R68 N; R50-53	T+; N R: 26/27/28-32-33-68-50/53 S: (1/2-)7-28-29-45-60-61	R32: C ≥ 1 % R33: C ≥ 0,1 %	
048-005-00-7	cadmiumhexafluorosilicate(2-); cadmium fluorosilica	241-084-0	17010-21-8	T; R23/25 R33 Xn; R68 N; R50-53	T; N R: 23/25-33-68-50/53 S: (1/2-)22-45-60-61	T; R23/25: C ≥ 10 % Xn; R20/22: 0,1 % ≤ C < 10 % R33: C ≥ 0,1 %	

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048-006-00-2	cadmium fluoride	232-222-0	7790-79-6	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23/25 N; R50-53	T+; N R: 45-46-60-61-25-26-48/23/25-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 % T; R25: C ≥ 10 % Xn; R22: 0,1 % ≤ C < 10 % T; R48/23/25: C ≥ 7 % Xn; R48/20/22: 0,1 % ≤ C < 7 %	E
048-007-00-8	cadmium iodide	232-223-6	7790-80-9	T; R23/25 R33 Xn; R68 N; R50-53	T; N R: 23/25-33-68-50/53 S: (1/2-)22-45-60-61	T; R23/25: C ≥ 10 % Xn; R20/22: 0,1 % ≤ C < 10 % R33: C ≥ 0,1 %	
048-008-00-3	cadmium chloride	233-296-7	10108-64-2	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23/25 N; R50-53	T+; N R: 45-46-60-61-25-26-48/23/25-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 % T; R25: C ≥ 10 % Xn; R22: 0,1 % ≤ C < 10 % T; R48/23/25: C ≥ 7 % Xn; R48/20/22: 0,1 % ≤ C < 7 %	E
048-009-00-9	cadmium sulphate	233-331-6	10124-36-4	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 T+; R26 T; R25-48/23/25 N; R50-53	T+; N R: 45-46-60-61-25-26-48/23/25-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 % T; R25: C ≥ 10 % Xn; R22: 0,1 % ≤ C < 10 % T; R48/23/25: C ≥ 7 % Xn; R48/20/22: 0,1 % ≤ C < 7 %	E

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048-010-00-4	cadmium sulphide	215-147-8	1306-23-6	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62-63 T; R48/23/25 Xn; R22 R53	T; R: 45-22-48/23/25-62-63-68-53 S: 53-45-61	Xn; R22: C ≥ 10 % T; R48/23/25: C ≥ 10 % Xn; R48/20/22: 0,1 % ≤ C < 10 %	E 1
048-011-00-X	cadmium (pyrophoric)	231-152-8	7440-43-9	F; R17 Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62-63 T+; R26 T; R48/23/25 N; R50-53	F; T+; N R: 45-17-26-48/23/25-62-63-68-50/53 S: 53-45-7/8-43-60-61		E
050-001-00-5	tin tetrachloride; stannic chloride	231-588-9	7646-78-8	C; R34 R52-53	C R: 34-52/53 S: (1/2-)7/8-26-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
050-002-00-0	cyhexatin (ISO); hydroxytricyclohexylstannane; tri(cyclohexyl)tin hydroxide	236-049-1	13121-70-5	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61		
050-003-00-6	fentin acetate (ISO); triphenyltin acetate	212-984-0	900-95-8	Carc. Cat. 3; R40 Repr. Cat. 3; R63 T+; R26 T; R24/25-48/23 Xi; R37/38-41 N; R50-53	T+; N R: 24/25-26-37/38-40-41-48/23-50/53-63 S: (1/2-)26-28-36/37/39-45-60-61		

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050-004-00-1	fentin hydroxide (ISO); triphenyltin hydroxide	200-990-6	76-87-9	Carc. Cat. 3; R40 Repr. Cat. 3; R63 T+; R26 T; R24/25-48/23 Xi; R37/38-41 N; R50-53	T+; N R: 24/25-26-37/38-40-41-48/23-50/53-63 S: (1/2-)26-28-36/37/39-45-60-61		
050-005-00-7	trimethyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)26-27-28-45-60-61	T+; R26/27/28: C ≥ 0,5 % T; R23/24/25: 0,1 % ≤ C < 0,5 % Xn; R20/21/22: 0,05 % ≤ C < 0,1 %	A 1
050-006-00-2	triethyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)26-27-28-45-60-61	T+; R26/27/28: C ≥ 0,5 % T; R23/24/25: 0,1 % ≤ C < 0,5 % Xn; R20/21/22: 0,05 % ≤ C < 0,1 %	A 1
050-007-00-8	tripropyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)26-27-28-45-60-61	T; R23/24/25: C ≥ 0,5 % Xn; R20/21/22: 0,1 % ≤ C < 0,5 %	A 1
050-008-00-3	tributyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	T; R25-48/23/25 Xn; R21 Xi; R36/38 N; R50-53	T; N R: 21-25-36/38-48/23/25-50/53 S: (1/2-)35-36/37/39-45-60-61	T; R25: C ≥ 1 % Xn; R22: 0,25 % ≤ C < 1 % T; R48/23/25: C ≥ 1 % Xn; R48/20/22: 0,25 % ≤ C < 1 % Xn; R21: C ≥ 1 % Xi; R36/38: C ≥ 1 %	A 1
050-009-00-9	fluorotripentylstannane; [1] hexapentyldistannoxane [2]	243-546-7 [1] 247-143-7 [2]	20153-49-5 [1] 25637-27-8 [2]	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)26-28-60-61	Xn; R20/21/22: C ≥ 1 %	1
050-010-00-4	fluorotrihexylstannane	243-547-2	20153-50-8	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)26-28-60-61	Xn; R20/21/22: C ≥ 1 %	1

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050-011-00-X	triphenyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)26-27-28-45-60-61	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,25 % ≤ C < 1 %	A 1
050-012-00-5	tetracyclohexylstannane; [1] chlorotricyclohexylstannane; [2] butyltricyclohexylstannane [3]	215-910-5 [1] 221-437-5 [2] 230-358-5 [3]	1449-55-4 [1] 3091-32-5 [2] 7067-44-9 [3]	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)26-28-60-61	Xn; R20/21/22: C ≥ 1 %	A 1
050-013-00-0	trioctyltin compounds, with the exception of those specified elsewhere in this Annex	—	—	Xi; R36/37/38 R53	Xi R: 36/37/38-53 S: (2-)61	Xi; R36/37/38: C ≥ 1 %	A 1
050-017-00-2	fenbutatin oxide (ISO); bis(tris(2-methyl-2-phenylpropyl)tin)oxide	236-407-7	13356-08-6	T+; R26 Xi; R36/38 N; R50-53	T+; N R: 26-36/38-50/53 S: (1/2-)28-36/37-45-60-61		
050-018-00-8	tin(II) methanesulphonate	401-640-7	53408-94-9	C; R34 Xn; R22 R43	C R: 22-34-43 S: (1/2-)22-26-36/37/39-45		
050-019-00-3	azocyclotin (ISO); 1-(tricyclohexylstannyl)-1*H*-1,2,4-triazole;	255-209-1	41083-11-8	T+; R26 T; R25 Xi; R37/38-41 N; R50-53	T+; N R: 25-26-37/38-41-50/53 S: (1/2-)26-28-36/37/39-38-45-60-61		
050-020-00-9	trioctylstannane	413-320-4	869-59-0	T; R48/25 Xi; R38 R53	T R: 38-48/25-53 S: (1/2-)23-36/37-45-61		
051-001-00-8	antimony trichloride	233-047-2	10025-91-9	C; R34 N; R51-53	C; N R: 34-51/53 S: (1/2-)26-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
051-002-00-3	antimony pentachloride	231-601-8	7647-18-9	C; R34 N; R51-53	C; N R: 34-51/53 S: (1/2-)26-45-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	

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051-003-00-9	antimony compounds, with the exception of the tetroxide (Sb_2O_4), pentoxide (Sb_2O_5), trisulphide (Sb_2S_3), pentasulphide (Sb_2S_5) and those specified elsewhere in this Annex	—	—	Xn; R20/22 N; R51-53	Xn; N R: 20/22-51/53 S: (2-)61	Xn; R20/22: C ≥ 0,25 %	A 1
051-004-00-4	antimony trifluoride	232-009-2	7783-56-4	T; R23/24/25 N; R51-53	T; N R: 23/24/25-51/53 S: (1/2-)7-26-45-61		
051-005-00-X	antimony trioxide	215-175-0	1309-64-4	Carc. Cat. 3; R40	Xn R: 40 S: (2-)22-36/37		
051-006-00-5	diphenyl(4-phenylthiophenyl)sulfonium hexafluoroantimonate	403-500-0	—	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
051-007-00-0	bis(4-dodecylphenyl)iodonium hexafluoroantimonate	404-420-9	71786-70-4	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
053-001-00-3	iodine	231-442-4	7553-56-2	Xn; R20/21 N; R50	Xn; N R: 20/21-50 S: (2-)23-25-61		
053-002-00-9	hydrogen iodide	233-109-9	10034-85-2	C; R35	C R: 35 S: (1/2-)9-26-36/37/39-45	C; R35: C ≥ 10 % C; R34: 0,2 % ≤ C < 10 % Xi; R36/37/38: 0,02 % ≤ C < 0,2 %	5
053-002-01-6	hydriodic acid ... %	—	—	C; R34	C R: 34 S: (1/2-)26-45	C; R34: C ≥ 25 % Xi; R36/38: 10 % ≤ C < 25 %	B
053-003-00-4	iodoxybenzene	—	696-33-3	E; R1 ⊗	E R: 1 S: (2-)35		
053-004-00-X	calcium iodoxybenzoate	—	—	E; R1 ⊗	E R: 1 S: (2-)35		C

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053-005-00-5	(4-(1-methylethyl)phenyl)-(4-methylphenyl)iodonium tetrakis (pentafluorophenyl)borate (1-)	422-960-3	178233-72-2	Xn; R21/22-48/22 N; R50-53	Xn; N R: 21/22-48/22-50/53 S: (2-)22-36/37-60-61		
056-001-00-1	barium peroxide	215-128-4	1304-29-6	O; R8 Xn; R20/22	O; Xn R: 8-20/22 S: (2-)13-27		
056-002-00-7	barium salts, with the exception of barium sulphate, salts of 1-azo-2-hydroxynaphthalenyl aryl sulphonic acid, and of salts specified elsewhere in this Annex	—	—	Xn; R20/22	Xn R: 20/22 S: (2-)28	Xn; R20/22: C ≥ 1 %	A 1
056-003-00-2	barium carbonate	208-167-3	513-77-9	Xn; R22	Xn R: 22 S: (2-)24/25		
056-004-00-8	barium chloride	233-788-1	10361-37-2	T; R25 Xn; R20	T R: 20-25 S: (1/2-)45		
072-001-00-4	hafnium tetra-*n*-butoxide	411-740-2	22411-22-9	Xi; R41 R43	Xi R: 41-43 S: (2-)24/25-26-37/39		
074-001-00-X	hexasodium tungstate hydrate	412-770-9	12141-67-2	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)22-26-39-61		
074-002-00-5	Reaction products of tungsten hexachloride with 2-methyl-propan-2-ol, nonylphenol and pentane-2,4-dione	408-250-6	—	F; R11 Xn; R20 C; R34 R43 N; R50-53	F; C; N R: 11-20-34-43-50/53 S: (1/2-)16-26-29-33-36/37/39-45-60-61		
076-001-00-5	osmium tetraoxide; osmic acid	244-058-7	20816-12-0	T+; R26/27/28 C; R34	T+ R: 26/27/28-34 S: (1/2-)7/9-26-45		
078-001-00-0	tetrachloroplatinates with the exception of those specified elsewhere in this Annex	—	—	T; R25 Xi; R41 R42/43	T R: 25-41-42/43 S: (2-)22-26-36/37/39-45		A

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078-002-00-6	diammonium tetrachloroplatinate	237-499-1	13820-41-2	T; R25 Xi; R38-41 R42/43	T R: 25-38-41-42/43 S: (2-)22-26-36/37/39-45		
078-003-00-1	disodium tetrachloroplatinate	233-051-4	10026-00-3	T; R25 Xi; R38-41 R42/43	T R: 25-38-41-42/43 S: (2-)22-26-36/37/39-45		
078-004-00-7	dipotassium tetrachloroplatinate	233-050-9	10025-99-7	T; R25 Xi; R38-41 R42/43	T R: 25-38-41-42/43 S: (2-)22-26-36/37/39-45		
078-005-00-2	hexachloroplatinates with the exception of those specified elsewhere in this Annex	—	—	T; R25 Xi; R41 R42/43	T R: 25-41-42/43 S: (1/2-)22-26-36/37/39-45		A
078-006-00-8	disodium hexachloroplatinate	240-983-5	16923-58-3	T; R25 Xi; R41 R42/43	T R: 25-41-42/43 S: (1/2-)22-26-36/37/39-45		
078-007-00-3	dipotassium hexachloroplatinate	240-979-3	16921-30-5	T; R25 Xi; R41 R42/43	T R: 25-41-42/43 S: (1/2-)22-26-36/37/39-45		
078-008-00-9	diammonium hexachloroplatinate	240-973-0	16919-58-7	T; R25 Xi; R41 R42/43	T R: 25-41-42/43 S: (1/2-)22-26-36/37/39-45		
078-009-00-4	hexachloroplatinic acid	241-010-7	16941-12-1	T; R25 C; R34 R42/43	T R: 25-34-42/43 S: (1/2-)22-26-36/37/39-45		
080-001-00-0	mercury	231-106-7	7439-97-6	T; R23 R33 N; R50-53	T; N R: 23-33-50/53 S: (1/2-)7-45-60-61		

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080-002-00-6	inorganic compounds of mercury with the exception of mercuric sulphide and those specified elsewhere in this Annex	—	—	T+; R26/27/28 R33 N; R50-53	T+; N R: 26/27/28-33-50/53 S: (1/2-)13-28-45-60-61	T+; R26/27/28: C ≥ 2 % T; R23/24/25: 0,5 % ≤ C < 2 % Xn; R20/21/22: 0,1 % ≤ C < 0,5 % R33: C ≥ 0,1 %	A 1
080-003-00-1	dimercury dichloride; mercurous chloride; calomel	233-307-5	10112-91-1	Xn; R22 Xi; R36/37/38 N; R50-53	Xn; N R: 22-36/37/38-50/53 S: (2-)13-24/25-46-60-61		
080-004-00-7	organic compounds of mercury with the exception of those specified elsewhere in this Annex	—	—	T+; R26/27/28 R33 N; R50-53	T+; N R: 26/27/28-33-50/53 S: (1/2-)13-28-36-45-60-61	T+; R26/27/28: C ≥ 2 % T; R23/24/25: 0,5 % ≤ C < 2 % Xn; R20/21/22: 0,05 % ≤ C < 0,5 % R33: C ≥ 0,05 %	A 1
080-005-00-2	mercury difulminate; mercuric fulminate; fulminate of mercury	211-057-8	628-86-4	E; R3 T; R23/24/25 R33 N; R50-53	E; T; N R: 3-23/24/25-33-50/53 S: (1/2-)3-45-60-61		
080-006-00-8	dimercury dicyanide oxide; mercuric oxycyanide	215-629-8	1335-31-5	E; R3 ⊗ T; R23/24/25 R33 N; R50-53	E; T; N R: 3-23/24/25-33-50/53 S: (1/2-)28-35-45-60-61		
080-007-00-3	dimethylmercury; [1] diethylmercury [2]	209-805-3 [1] 211-000-7 [2]	593-74-8 [1] 627-44-1 [2]	T+; R26/27/28 R33 N; R50-53	T+; N R: 26/27/28-33-50/53 S: (1/2-)13-28-36-45-60-61	T+; R26/27/28: C ≥ 0,5 % T; R23/24/25: 0,1 % ≤ C < 0,5 % Xn; R20/21/22: 0,05 % ≤ C < 0,1 % R33: C ≥ 0,05 %	1
080-008-00-9	phenylmercury nitrate; [1] phenylmercury hydroxide; [2] basic phenylmercury nitrate [3]	200-242-9 [1] 202-866-7 [2] - [3]	55-68-5 [1] 100-57-2 [2] 8003-05-2 [3]	T; R25-48/24/25 C; R34 N; R50-53	T; N R: 25-34-48/24/25-50/53 S: (1/2-)23-24/25-37-45-60-61		

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080-009-00-4	2-methoxyethylmercury chloride	204-659-7	123-88-6	T; R25-48/25 C; R34 N; R50-53	T; N R: 25-34-48/25-50/53 S: (1/2-)36/37/39-45-60-61		
080-010-00-X	mercury dichloride; mercuric chloride	231-299-8	7487-94-7	T+; R28 T; R48/24/25 C; R34 N; R50-53	T+; N R: 28-34-48/24/25-50/53 S: (1/2-)36/37/39-45-60-61		
080-011-00-5	phenylmercury acetate	200-532-5	62-38-4	T; R25-48/24/25 C; R34 N; R50-53	T; N R: 25-34-48/24/25-50/53 S: (1/2-)23-24/25-37-45-60-61		
081-001-00-3	thallium	231-138-1	7440-28-0	T+; R26/28 R33 R53	T+ R: 26/28-33-53 S: (1/2-)13-28-45-61		
081-002-00-9	thallium compounds, with the exception of those specified elsewhere in this Annex	—	—	T+; R26/28 R33 N; R51-53	T+; N R: 26/28-33-51/53 S: (1/2-)13-28-45-61		A
081-003-00-4	dithallium sulphate; thallic sulphate	231-201-3	7446-18-6	T+; R28 T; R48/25 Xi; R38 N; R51-53	T+; N R: 28-38-48/25-51/53 S: (1/2-)13-36/37-45-61		
082-001-00-6	lead compounds with the exception of those specified elsewhere in this Annex	—	—	Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R20/22 R33 N; R50-53	T; N R: 61-20/22-33-62-50/53 S: 53-45-60-61	Repr. Cat. 3; R62: C ≥ 2,5 % Xn; R20/22: C ≥ 1 % R33: C ≥ 0,5 %	AE 1

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082-002-00-1	lead alkyls	—	—	Repr. Cat. 1; R61 Repr. Cat. 3; R62 T+; R26/27/28 R33 N; R50-53	T+; N R: 61-26/27/28-33-62-50/53 S: 53-45-60-61	Repr. Cat. 1; R61: C ≥ 0,1 % T+; R26/27/28: C ≥ 0,25 % T; R23/24/25: 0,1 % ≤ C < 0,25 % Xn; R20/21/22: 0,05 % ≤ C < 0,1 % R33: C ≥ 0,05 %	AE 1
082-003-00-7	lead diazide; lead azide	236-542-1	13424-46-9	E; R3 Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R20/22 R33 N; R50-53	E; T; N R: 61-3-20/22-33-50/53-62 S: 53-45-60-61		E 1
082-004-00-2	lead chromate	231-846-0	7758-97-6	Carc. Cat. 3; R40 Repr. Cat. 1; R61 Repr. Cat. 3; R62 R33 N; R50-53	T; N R: 61-33-40-50/53-62 S: 53-45-60-61		1
082-005-00-8	lead di(acetate)	206-104-4	301-04-2	Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R48/22 R33 N; R50-53	T; N R: 61-33-48/22-50/53-62 S: 53-45-60-61		E 1
082-006-00-3	trilead bis(orthophosphate)	231-205-5	7446-27-7	Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R48/22 R33 N; R50-53	T; N R: 61-33-48/22-50/53-62 S: 53-45-60-61		E 1

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082-007-00-9	lead acetate, basic;	215-630-3	1335-32-6	Carc. Cat. 3; R40 Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R48/22 R33 N; R50-53	T; N R: 61-33-40-48/22-50/53-62 S: 53-45-60-61		E 1
082-008-00-4	lead(II) methanesulphonate	401-750-5	17570-76-2	Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R20/22-48/20/22 Xi; R38-41 N; R58 R33	T; N R: 61-62-20/22-33-38-41-48/20/22-58 S: 53-45-57-61		E 1
082-009-00-X	Lead sulfochromate yellow; C.I. Pigment Yellow 34; [This substance is identified in the Colour Index by Colour Index Constitution Number, C.I. 77603.]	215-693-7	1344-37-2	Carc. Cat. 3; R40 Repr. Cat. 1; R61 Repr. Cat. 3; R62 R33 N; R50-53	T; N R: 61-33-40-50/53-62 S: 53-45-60-61		1
082-010-00-5	Lead chromate molybdate sulfate red; C.I. Pigment Red 104; [This substance is identified in the Colour Index by Colour Index Constitution Number, C.I. 77605.]	235-759-9	12656-85-8	Carc. Cat. 3; R40 Repr. Cat. 1; R61 Repr. Cat. 3; R62 R33 N; R50-53	T; N R: 61-33-40-50/53-62 S: 53-45-60-61		1

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082-011-00-0	lead hydrogen arsenate	232-064-2	7784-40-9	Carc. Cat. 1; R45 Repr. Cat. 1; R61 Repr. Cat. 3; R62 T; R23/25 R33 N; R50-53	T; N R: 45-61-23/25-33-50/53-62 S: 53-45-60-61		E 1
092-001-00-8	uranium	231-170-6	7440-61-1	T+; R26/28 R33 R53	T+ R: 26/28-33-53 S: (1/2-)20/21-45-61		
092-002-00-3	uranium compounds	—	—	T+; R26/28 R33 N; R51-53	T+; N R: 26/28-33-51/53 S: (1/2-)20/21-45-61		A
601-001-00-4	methane	200-812-7	74-82-8	F+; R12	F+ R: 12 S: (2-)9-16-33		
601-002-00-X	ethane	200-814-8	74-84-0	F+; R12	F+ R: 12 S: (2-)9-16-33		
601-003-00-5	propane	200-827-9	74-98-6	F+; R12	F+ R: 12 S: (2-)9-16		
601-004-00-0	butane; [1] and isobutane [2]	203-448-7 [1] 200-857-2 [2]	106-97-8 [1] 75-28-5 [2]	F+; R12	F+ R: 12 S: (2-)9-16		C
601-004-01-8	butane (containing ≥ 0,1 % butadiene (203-450-8)); [1] isobutane (containing ≥ 0,1 % butadiene (203-450-8)) [2]	203-448-7 [1] 200-857-2 [2]	106-97-8 [1] 75-28-5 [2]	F+; R12 Carc. Cat. 1; R45 Muta. Cat. 2; R46	F+; T R: 45-46-12 S: 53-45		CS
601-005-00-6	2,2-dimethylpropane; neopentane	207-343-7	463-82-1	F+; R12 N; R51-53	F+; N R: 12-51/53 S: (2-)9-16-33-61		

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601-006-00-1	pentane; [1] isopentane; 2-methylbutane [2]	203-692-4 [1] 201-142-8 [2]	109-66-0 [1] 78-78-4 [2]	F+; R12 Xn; R65 R66 R67 N; R51-53	F+; Xn; N R: 12-51/53-65-66-67 S: (2-)9-16-29-33-61-62		C
601-007-00-7	hexane, reaction mass of isomers (containing < 5 % *n*-hexane (203-777-6))	—	—	F; R11 Xn; R65 Xi; R38 R67 N; R51-53	F; Xn; N R: 11-38-51/53-65-67 S: (2-)9-16-29-33-61-62		C
601-008-00-2	heptane [and isomers] [1]	205-563-8 [1] 203-548-0 [2] 207-346-3 [3] 209-230-8 [4] 209-280-0 [5] 209-643-3 [6] 209-680-5 [7] 209-730-6 [8] 210-529-0 [9] 250-610-8 [10]	142-82-5 [1] 108-08-7 [2] 464-06-2 [3] 562-49-2 [4] 565-59-3 [5] 589-34-4 [6] 590-35-2 [7] 591-76-4 [8] 617-78-7 [9] 31394-54-4 [10]	F; R11 Xn; R65 Xi; R38 R67 N; R50-53	F; Xn; N R: 11-38-50/53-65-67 S: (2-)9-16-29-33-60-61-62		C

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601-009-00-8	octane [and isomers] [1]	203-892-1 [1] 208-759-1 [2] 209-207-2 [3] 209-243-9 [4] 209-266-4 [5] 209-292-6 [6] 209-504-7 [7] 209-547-1 [8] 209-649-6 [9] 209-650-1 [10] 209-660-6 [11] 209-689-4 [12]	111-65-9 [1] 540-84-1 [2] 560-21-4 [3] 563-16-6 [4] 564-02-3 [5] 565-75-3 [6] 583-48-2 [7] 584-94-1 [8] 589-43-5 [9] 589-53-7 [10] 589-81-1 [11] 590-73-8 [12] 592-13-2 [13] 592-27-8 [14] 594-82-1 [15] 609-26-7 [16] 619-99-8 [17] 1067-08-9 [18] 26635-64-3 [19]	F; R11 Xn; R65 Xi; R38 R67 N; R50-53	F; Xn; N R: 11-38-50/53-65-67 S: (2-)9-16-29-33-60-61-62		C
		209-745-8 [13] 209-747-9 [14] 209-855-6 [15] 210-187-2 [16] 210-621-0 [17] 213-923-0 [18] 247-861-0 [19]					

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601-010-00-3	ethylene	200-815-3	74-85-1	F+; R12 R67	F+ R: 12-67 S: (2-)9-16-33-45		
601-011-00-9	propene; propylene	204-062-1	115-07-1	F+; R12	F+ R: 12 S: (2-)9-16-33		
601-012-00-4	but-1-ene; [1] butene, mixed-1-and-2-isomers; [2] 2-methylpropene; [3] (*Z*)-but-2-ene; [4] (*E*)-but-2-ene [5]	203-449-2 [1] 203-452-9 [2] 204-066-3 [3] 209-673-7 [4] 210-855-3 [5]	106-98-9 [1] 107-01-7 [2] 115-11-7 [3] 590-18-1 [4] 624-64-6 [5]	F+; R12	F+ R: 12 S: (2-)9-16-33		C
601-013-00-X	1,3-butadiene; buta-1,3-diene	203-450-8	106-99-0	F+; R12 Carc. Cat. 1; R45 Muta. Cat. 2; R46	F+; T R: 45-46-12 S: 53-45		D
601-014-00-5	isoprene (stabilised) 2-methyl-1,3-butadiene	201-143-3	78-79-5	F+; R12 Carc. Cat. 2; R45 Muta. Cat. 3; R68 R52-53	F+; T R: 45-12-68-52/53 S: 53-45-61		D
601-015-00-0	acetylene; ethyne	200-816-9	74-86-2	R5 R6 F+; R12	F+ R: 5-6-12 S: (2-)9-16-33		
601-016-00-6	cyclopropane	200-847-8	75-19-4	F+; R12	F+ R: 12 S: (2-)9-16-33		
601-017-00-1	cyclohexane	203-806-2	110-82-7	F; R11 Xn; R65 Xi; R38 R67 N; R50-53	F; Xn; N R: 11-38-65-67-50/53 S: (2-)9-16-25-33-60-61-62		

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601-018-00-7	methylcyclohexane	203-624-3	108-87-2	F; R11 Xn; R65 Xi; R38 R67 N; R51-53	F; Xn; N R: 11-38-51/53-65-67 S: (2-)9-16-33-61-62		
601-019-00-2	1,4-dimethylcyclohexane	209-663-2	589-90-2	F; R11 Xn; R65 Xi; R38 R67 N; R51-53	F; Xn; N R: 11-38-51/53-65-67 S: (2-)9-16-33-61-62		
601-020-00-8	benzene	200-753-7	71-43-2	F; R11 Carc. Cat. 1; R45 Muta. Cat. 2; R46 T; R48/23/24/25 Xn; R65 Xi; R36/38	F; T R: 45-46-11-36/38-48/23/24/25-65 S: 53-45		E
601-021-00-3	toluene	203-625-9	108-88-3	F; R11 Repr. Cat. 3; R63 Xn; R48/20-65 Xi; R38 R67	F; Xn R: 11-38-48/20-63-65-67 S: (2-)36/37-46-62		
601-022-00-9	*o*-xylene; [1] *p*-xylene; [2] *m*-xylene; [3] xylene [4]	202-422-2 [1] 203-396-5 [2] 203-576-3 [3] 215-535-7 [4]	95-47-6 [1] 106-42-3 [2] 108-38-3 [3] 1330-20-7 [4]	R10 Xn; R20/21 Xi; R38	Xn R: 10-20/21-38 S: (2-)25	Xn; R20/21: C ≥ 12,5 %	C
601-023-00-4	ethylbenzene	202-849-4	100-41-4	F; R11 Xn; R20	F; Xn R: 11-20 S: (2-)16-24/25-29		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
601-024-00-X	cumene; [1] propylbenzene [2]	202-704-5 [1] 203-132-9 [2]	98-82-8 [1] 103-65-1 [2]	R10 Xn; R65 Xi; R37 N; R51-53	Xn; N R: 10-37-51/53-65 S: (2-)24-37-61-62		C
601-025-00-5	mesitylene; 1,3,5-trimethylbenzene	203-604-4	108-67-8	R10 Xi; R37 N; R51-53	Xi; N R: 10-37-51/53 S: (2-)61	Xi; R37: C ≥ 25 %	
601-026-00-0	styrene	202-851-5	100-42-5	R10 Xn; R20 Xi; R36/38	Xn R: 10-20-36/38 S: (2-)23	Xn; R20: C ≥ 12,5 % Xi; R36/38: C ≥ 12,5 %	D
601-027-00-6	2-phenylpropene; α-methylstyrene	202-705-0	98-83-9	R10 Xi; R36/37 N; R51-53	Xi; N R: 10-36/37-51/53 S: (2-)61	Xi; R36/37: C ≥ 25 %	
601-028-00-1	2-methylstyrene; 2-vinyltoluene	210-256-7	611-15-4	Xn; R20 N; R51-53	Xn; N R: 20-51/53 S: (2-)24-61		
601-029-00-7	dipentene; limonene; [1] (*R*)-*p*-mentha-1,8-diene; d-limonene; [2] (*S*)-*p*-mentha-1,8-diene; l-limonene; [3] *trans*-1-methyl-4-(1-methylvinyl)cyclohexene; [4] (±)-1-methyl-4-(1-methylvinyl)cyclohexene [5]	205-341-0 [1] 227-813-5 [2] 227-815-6 [3] 229-977-3 [4] 231-732-0 [5]	138-86-3 [1] 5989-27-5 [2] 5989-54-8 [3] 6876-12-6 [4] 7705-14-8 [5]	R10 Xi; R38 R43 N; R50-53	Xi; N R: 10-38-43-50/53 S: (2-)24-37-60-61		C
601-030-00-2	cyclopentane	206-016-6	287-92-3	F; R11 R52-53	F R: 11-52/53 S: (2-)9-16-29-33-61		
601-031-00-8	2,4,4-trimethylpent-1-ene	203-486-4	107-39-1	F; R11 N; R51-53	F; N R: 11-51/53 S: (2-)9-16-29-33-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
601-032-00-3	benzo[*a*]pyrene; benzo[*def*]chrysene	200-028-5	50-32-8	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 R43 N; R50-53	T; N R: 45-46-60-61-43-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 %	
601-033-00-9	benz[*a*]anthracene	200-280-6	56-55-3	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
601-034-00-4	benz[*e*]acephenanthrylene	205-911-9	205-99-2	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
601-035-00-X	benzo[*j*]fluoranthene	205-910-3	205-82-3	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
601-036-00-5	benzo[*k*]fluoranthene	205-916-6	207-08-9	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
601-037-00-0	n-hexane	203-777-6	110-54-3	F; R11 Repr. Cat. 3; R62 Xn; R65-48/20 Xi; R38 R67 N; R51-53	F; Xn; N R: 11-38-48/20-62-65-67-51/53 S: (2-)9-16-29-33-36/37-61-62	Xn; R48/20: C ≥ 5 %	
601-041-00-2	dibenz[*a*,*h*]anthracene	200-181-8	53-70-3	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 %	
601-042-00-8	biphenyl; diphenyl	202-163-5	92-52-4	Xi; R36/37/38 N; R50-53	Xi; N R: 36/37/38-50/53 S: (2-)23-60-61		

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601-043-00-3	1,2,4-trimethylbenzene	202-436-9	95-63-6	R10 Xn; R20 Xi; R36/37/38 N; R51-53	Xn; N R: 10-20-36/37/38-51/53 S: (2-)26-61		
601-044-00-9	3a,4,7,7a-tetrahydro-4,7-methanoindene	201-052-9	77-73-6	F; R11 Xn; R20/22 Xi; R36/37/38 N; R51-53	F; Xn; N R: 11-20/22-36/37/38-51/53 S: (2-)36/37-61		
601-045-00-4	1,2,3,4-tetrahydronaphthalene	204-340-2	119-64-2	R19 Xi; R36/38 N; R51-53	Xi; N R: 19-36/38-51/53 S: (2-)26-28-61		
601-046-00-X	7-methylocta-1,6-diene	404-210-7	42152-47-6	R10 N; R50-53	N R: 10-50/53 S: (2-)60-61		
601-047-00-5	*m*-mentha-1,3(8)-diene	404-150-1	17092-80-7	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
601-048-00-0	chrysene	205-923-4	218-01-9	Carc. Cat. 2; R45 Muta. Cat. 3; R68 N; R50-53	T; N R: 45-68-50/53 S: 53-45-60-61		
601-049-00-6	benzo[*e*]pyrene	205-892-7	192-97-2	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
601-051-00-7	4-phenylbut-1-ene	405-980-7	768-56-9	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
601-052-00-2	naphthalene	202-049-5	91-20-3	Carc. Cat. 3; R40 Xn; R22 N; R50-53	Xn; N R: 22-40-50/53 S: (2-)36/37-46-60-61		

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601-053-00-8	nonylphenol; [1] 4-nonylphenol, branched [2]	246-672-0 [1] 284-325-5 [2]	25154-52-3 [1] 84852-15-3 [2]	Repr. Cat. 3; R62-63 Xn; R22 C; R34 N; R50-53	C; N R: 22-34-62-63-50/53 S: (1/2-)26-36/37/39-45-46-60-61		
601-054-00-3	reaction mass of isomers of: dibenzylbenzene; dibenzyl(methyl)benzene; dibenzyl(dimethyl)benzene; dibenzyl(trimethyl)benzene	405-570-8	—	N; R50-53	N R: 50/53 S: 60-61		
601-055-00-9	reaction mass of isomers of: mono-(2-tetradecyl)naphthalenes; di-(2-tetradecyl)naphthalenes; tri-(2-tetradecyl)naphthalenes	410-190-0	132983-41-6	Xi; R36 R53	Xi R: 36-53 S: (2-)26-61		
601-056-00-4	reaction mass of isomers of: methyldiphenylmethane; dimethyldiphenylmethane	405-470-4	73807-39-3	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)37-60-61		
601-057-00-X	*N*-dodecyl-[3-(4-(dimethylamino)benzamido)-propyl] dimethylammonium tosylate	421-130-8	156679-41-3	Xi; R41 R43 N; R50-53	Xi; N R: 41-43-50/53 S: (2-)24-26-37/39-60-61		
601-058-00-5	di-L-para-menthene	417-870-6	83648-84-4	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)23-24-37-60-61		
601-059-00-0	methyl 2-benzylidene-3-oxobutyrate	420-940-9	15768-07-7	Xi; R36/38 N; R51-53	Xi; N R: 36/38-51/53 S: (2-)26-37/39-61		
601-060-00-6	1,2-bis[4-fluoro-6-{4-sulfo-5-(2-(4-sulfonaphtalene-3-ylazo)-1-hydroxy-3,6-disulfo-8-aminonaphthalene-7-ylazo)phenylamino}-1,3,5-triazin-2ylamino]ethane; x-sodium, y-potassium salts x = 7,755 y = 0,245	417-610-1	155522-09-1	R43	Xi R: 43 S: (2-)22-24-37		
601-061-00-1	(ethyl-1,2-ethanediyl)[-2-[[[(2-hydroxyethyl)methylamino] acetyl]-propyl]ω-(nonylphenoxy)poly]oxy-(methyl-1,2-ethanediyl)	418-960-8	—	C; R34 R43 N; R51-53	C; N R: 34-43-51/53 S: (1/2-)26-28-36/37/39-45-61		

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601-062-00-7	reaction mass of: branched triacontane; branched dotriacontane; branched tetratriacontane; branched hexatriacontane	417-030-9	151006-59-6	R53	R: 53 S: 61		
601-063-00-2	reaction mass of isomers of branched tetracosane	417-060-2	151006-61-0	Xn; R20 R53	Xn R: 20-53 S: (2-)61		
601-064-00-8	branched hexatriacontane	417-070-7	151006-62-1	R53	R: 53 S: 61		
601-065-00-3	reaction mass of: (1'-α,3'-α,6'-α-2,2,3',7',7'-pentamethylspiro (1,3-dioxane-5,2'-norcarane); (1'α,3'β,6'α)-2,2,3',7',7'-pentamethylspiro(1,3-dioxane-5,2'-norcarane)	416-930-9	—	Xn; R48/22 Xi; R41 N; R51-53	Xn; N R: 41-48/22-51/53 S: (2-)22-26-37/39-61		
601-066-00-9	1-(4-(*trans*-4-heptylcyclohexyl)phenyl)ethanone	426-820-2	78531-60-9	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
601-067-00-4	triethyl arsenate	427-700-2	15606-95-8	Carc. Cat. 1; R45 T; R23/25 N; R50-53	T; N R: 45-23/25-50/53 S: 53-45-60-61		E
601-068-00-X	1,2-diacetoxybut-3-ene	421-720-5	18085-02-4	Xn; R22	Xn R: 22 S: (2-)		
601-069-00-5	2-ethyl-1-(2-(1,3-dioxanyl)ethyl)-pyridinium bromide	422-680-1	287933-44-2	R52-53	R: 52/53 S: 61		
601-071-00-6	1-dimethoxymethyl-2-nitro-benzene	423-830-9	20627-73-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
601-073-00-7	1-bromo-3,5-difluorobenzene	416-710-2	461-96-1	R10 Xn; R22-48/22 Xi; R38 R43 N; R50-53	Xn; N R: 10-22-38-43-48/22-50/53 S: (2-)24-36/37-60-61		

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601-074-00-2	reaction mass of: 4-(2,2,3-trimethylcyclopent-3-en-1-yl)-1-methyl-2-oxabicyclo[2.2.2]octane; 1-(2,2,3-trimethylcyclopent-3-en-1-yl)-5-methyl-6-oxabicyclo[3.2.1]octane; spiro[cyclohex-3-en-1-yl-[(4,5,6,6a-tetrahydro-3,6',6',6'a-tetramethyl)-1,3'(3'a*H*)-[2*H*]cyclopenta[*b*]furan]; spiro[cyclohex-3-en-1-yl-[4,5,6,6a-tetrahydro-4,6',6',6'a-tetramethyl)-1,3'(3'a*H*)-[2*H*]cyclopenta[*b*]]furan]	422-040-1	—	Xi; R36/38 N; R51-53	Xi; N R: 36/38-51/53 S: (2-)26-37-61		
602-001-00-7	chloromethane; methyl chloride	200-817-4	74-87-3	F+; R12 Carc. Cat. 3; R40 Xn; R48/20	F+; Xn R: 12-40-48/20 S: (2-)9-16-33		
602-002-00-2	bromomethane; methylbromide	200-813-2	74-83-9	Muta. Cat. 3; R68 T; R23/25 Xn; R48/20 Xi; R36/37/38 N; R50 N; R59	T; N R: 23/25-36/37/38-48/20-50-59-68 S: (1/2-)15-27-36/39-38-45-59-61		
602-003-00-8	dibromomethane	200-824-2	74-95-3	Xn; R20 R52-53	Xn R: 20-52/53 S: (2-)24-61	Xn; R20: C ≥ 12,5 %	
602-004-00-3	dichloromethane; methylene chloride	200-838-9	75-09-2	Carc. Cat. 3; R40	Xn R: 40 S: (2-)23-24/25-36/37		
602-005-00-9	methyl iodide; iodomethane	200-819-5	74-88-4	Carc. Cat. 3; R40 Xn; R21 T; R23/25 Xi; R37/38	T R: 21-23/25-37/38-40 S: (1/2-)36/37-38-45		
602-006-00-4	trichloromethane; chloroform	200-663-8	67-66-3	Xn; R22-48/20/22 Xi; R38 Carc. Cat. 3; R40	Xn R: 22-38-40-48/20/22 S: (2-)36/37	Xn; R22: C ≥ 5 % Xn; R48/20/22: C ≥ 5 %	

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602-007-00-X	bromoform; tribromomethane	200-854-6	75-25-2	T; R23 Xi; R36/38 N; R51-53	T; N R: 23-36/38-51/53 S: (1/2-)28-45-61		
602-008-00-5	carbon tetrachloride; tetrachloromethane	200-262-8	56-23-5	Carc. Cat. 3; R40 T; R23/24/25-48/23 N; R59 R52-53	T; N R: 23/24/25-40-48/23-59-52/53 S: (1/2-)23-36/37-45-59-61	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,2 % ≤ C < 1 % T; R48/23: C ≥ 1 % Xn; R48/20: 0,2 % ≤ C < 1 %	
602-009-00-0	chloroethane	200-830-5	75-00-3	F+; R12 Carc. Cat. 3; R40 R52-53	F+; Xn R: 12-40-52/53 S: (2-)9-16-33-36/37-61		
602-010-00-6	1,2-dibromoethane	203-444-5	106-93-4	Carc. Cat. 2; R45 T; R23/24/25 Xi; R36/37/38 N; R51-53	T; N R: 45-23/24/25-36/37/38-51/53 S: 53-45-61	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,1 % ≤ C < 1 %	E
602-011-00-1	1,1-dichloroethane	200-863-5	75-34-3	F; R11 Xn; R22 Xi; R36/37 R52-53	F; Xn R: 11-22-36/37-52/53 S: (2-)16-23-61	Xn; R22: C ≥ 12,5 %	
602-012-00-7	1,2-dichloroethane; ethylene dichloride	203-458-1	107-06-2	F; R11 Carc. Cat. 2; R45 Xn; R22 Xi; R36/37/38	F; T R: 45-11-22-36/37/38 S: 53-45		E
602-013-00-2	1,1,1-trichloroethane; methyl chloroform	200-756-3	71-55-6	Xn; R20 N; R59	Xn; N R: 20-59 S: (2-)24/25-59-61		F
602-014-00-8	1,1,2-trichloroethane	201-166-9	79-00-5	Carc. Cat. 3; R40 Xn; R20/21/22 R66	Xn R: 20/21/22-40-66 S: (2-)9-36/37-46	Xn; R20/21/22: C ≥ 5 %	

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602-015-00-3	1,1,2,2-tetrachloroethane	201-197-8	79-34-5	T+; R26/27 N; R51-53	T+; N R: 26/27-51/53 S: (1/2-)38-45-61		
602-016-00-9	1,1,2,2-tetrabromoethane	201-191-5	79-27-6	T+; R26 Xi; R36 R52-53	T+ R: 26-36-52/53 S: (1/2-)24-27-45-61		
602-017-00-4	pentachloroethane	200-925-1	76-01-7	Carc. Cat. 3; R40 T; R48/23 N; R51-53	T; N R: 40-48/23-51/53 S: (1/2-)23-36/37-45-61	T; R48/23: C ≥ 1 % Xn; R48/20: 0,2 % ≤ C < 1 %	
602-018-00-X	1-chloropropane; [1] 2-chloropropane [2]	208-749-7 [1] 200-858-8 [2]	540-54-5 [1] 75-29-6 [2]	F; R11 Xn; R20/21/22	F; Xn R: 11-20/21/22 S: (2-)9-29		C
602-019-00-5	1-bromopropane; n-propyl bromide	203-445-0	106-94-5	F; R11 Repr. Cat. 2; R60 Repr. Cat. 3; R63 Xn; R48/20 Xi; R36/37/38 R67	F; T R: 60-11-36/37/38-48/20-63-67 S: 53-45		E
602-020-00-0	1,2-dichloropropane; propylene dichloride	201-152-2	78-87-5	F; R11 Xn; R20/22	F; Xn R: 11-20/22 S: (2-)16-24		
602-021-00-6	1,2-dibromo-3-chloropropane	202-479-3	96-12-8	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 1; R60 T; R25 Xn; R48/20/22 R52-53	T R: 45-46-60-25-48/20/22-52/53 S: 53-45-61		E

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602-022-00-1	1-chloropentane; [1] 2-chloropentane; [2] 3-chloropentane [3]	208-846-4 [1] 210-885-7 [2] 210-467-4 [3]	543-59-9 [1] 625-29-6 [2] 616-20-6 [3]	F; R11 Xn; R20/21/22	F; Xn R: 11-20/21/22 S: (2-)9-29		C
602-023-00-7	vinyl chloride; chloroethylene	200-831-0	75-01-4	F+; R12 Carc. Cat. 1; R45	F+; T R: 45-12 S: 53-45		D
602-024-00-2	bromoethylene	209-800-6	593-60-2	F+; R12 Carc. Cat. 2; R45	F+; T R: 45-12 S: 53-45		
602-025-00-8	1,1-dichloroethylene; vinylidene chloride	200-864-0	75-35-4	F+; R12 Carc. Cat. 3; R40 Xn; R20	F+; Xn R: 12-20-40 S: (2-)7-16-29-36/37-46	Xn; R20: C ≥ 12,5 %	D
602-026-00-3	1,2-dichloroethylene; [1] *cis*-dichloroethylene; [2] *trans*-dichloroethylene [3]	208-750-2 [1] 205-859-7 [2] 205-860-2 [3]	540-59-0 [1] 156-59-2 [2] 156-60-5 [3]	F; R11 Xn; R20 R52-53	F; Xn R: 11-20-52/53 S: (2-)7-16-29-61	Xn; R20: C ≥ 12,5 %	C
602-027-00-9	trichloroethylene; trichloroethene	201-167-4	79-01-6	Carc. Cat. 2; R45 Muta. Cat. 3; R68 R67 Xi; R36/38 R52-53	T R: 45-36/38-52/53-67 S: 53-45-61		
602-028-00-4	tetrachloroethylene	204-825-9	127-18-4	Carc. Cat. 3; R40 N; R51-53	Xn; N R: 40-51/53 S: (2-)23-36/37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
602-029-00-X	3-chloropropene; allyl chloride	203-457-6	107-05-1	F; R11 Carc. Cat. 3; R40 Muta. Cat. 3; R68 Xn; R20/21/22-48/20 Xi; R36/37/38 N; R50	F; Xn; N R: 11-20/21/22-36/37/38-40-48/20-68-50 S: (2-)16-25-26-36/37-46-61		D
602-030-00-5	1,3-dichloropropene; [1] (Z)-1,3-dichloropropene [2]	208-826-5 [1] 233-195-8 [2]	542-75-6 [1] 10061-01-5 [2]	R10 T; R25 Xn; R20/21 Xi; R36/37/38 R43 N; R50-53	T; N R: 10-20/21-25-36/37/38-43-50/53 S: (1/2-)36/37-45-60-61		DC
602-031-00-0	1,1-dichloropropene	209-253-3	563-58-6	F; R11 T; R25 R52-53	F; T R: 11-25-52/53 S: (1/2-)16-29-33-45-61		
602-032-00-6	3-chloro-2-methylpropene	209-251-2	563-47-3	F; R11 Xn; R20/22 C; R34 R43 N; R51-53	F; C; N R: 11-20/22-34-43-51/53 S: (2-)9-16-26-29-36/37/39-45-61		
602-033-00-1	chlorobenzene	203-628-5	108-90-7	R10 Xn; R20 N; R51-53	Xn; N R: 10-20-51/53 S: (2-)24/25-61	Xn; R20: C ≥ 5 %	
602-034-00-7	1,2-dichlorobenzene; *o*-dichlorobenzene	202-425-9	95-50-1	Xn; R22 Xi; R36/37/38 N; R50-53	Xn; N R: 22-36/37/38-50/53 S: (2-)23-60-61	Xn; R22: C ≥ 5 %	
602-035-00-2	1,4-dichlorobenzene; *p*-dichlorobenzene	203-400-5	106-46-7	Carc. Cat. 3; R40 Xi; R36 N; R50-53	Xn; N R: 36-40-50/53 S: (2-)36/37-46-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
602-036-00-8	chloroprene (stabilised); 2-chlorobuta-1,3-diene (stabilised)	204-818-0	126-99-8	F; R11 Carc. Cat. 2; R45 Xn; R20/22-48/20 Xi; R36/37/38	F; T R: 45-11-20/22-36/37/38-48/20 S: 53-45		D E
602-037-00-3	α-chlorotoluene; benzyl chloride	202-853-6	100-44-7	Carc. Cat. 2; R45 T; R23 Xn; R22-48/22 Xi; R37/38-41	T R: 45-22-23-37/38-41-48/22 S: 53-45		E
602-038-00-9	α,α,α-trichlorotoluene; benzotrichloride	202-634-5	98-07-7	Carc. Cat. 2; R45 T; R23 Xn; R22 Xi; R37/38-41	T R: 45-22-23-37/38-41 S: 53-45		E
602-039-00-4	polychlorobiphenyls; PCB	215-648-1	1336-36-3	R33 N; R50-53	N R: 33-50/53 S: (2-)-60-61	R33: C ≥ 0,005 %	C
602-040-00-X	2-chlorotoluene; [1] 3-chlorotoluene; [2] 4-chlorotoluene; [3] chlorotoluene [4]	202-424-3 [1] 203-580-5 [2] 203-397-0 [3] 246-698-2 [4]	95-49-8 [1] 108-41-8 [2] 106-43-4 [3] 25168-05-2 [4]	Xn; R20 N; R51-53	Xn; N R: 20-51/53 S: (2-)24/25-61		C
602-041-00-5	penthachloronaphthalene	215-320-8	1321-64-8	Xn; R21/22 Xi; R36/38 N; R50-53	Xn; N R: 21/22-36/38-50/53 S: (2-)-60-61		C
602-042-00-0	1,2,3,4,5,6-hexachlorcyclohexanes with the exception of those specified elsewhere in this Annex	—	—	Carc. Cat. 3; R40 T; R25 Xn; R21 N; R50-53	T; N R: 21-25-40-50/53 S: (1/2-)22-36/37-45-60-61		A C

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
602-043-00-6	lindane (ISO) γ-HCH or γ-BHC; γ-1,2,3,4,5,6-hexachlorocyclohexane;	200-401-2	58-89-9	T; R25 Xn; R20/21-48/22 R64 N; R50-53	T; N R: 20/21-25-48/22-64-50/53 S: (1/2-)36/37-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
602-044-00-1	camphechlor (ISO) toxaphene;	232-283-3	8001-35-2	Carc. Cat. 3; R40 T; R25 Xn; R21 Xi; R37/38 N; R50-53	T; N R: 21-25-37/38-40-50/53 S: (1/2-)36/37-45-60-61		
602-045-00-7	DDT (ISO); clofenotane (INN); dicophane; 1,1,1-trichloro-2,2-bis(4-chlorophenyl)ethane; dichlorodiphenyltrichloroethane	200-024-3	50-29-3	T; R25-48/25 Carc. Cat. 3; R40 N; R50-53	T; N R: 25-40-48/25-50/53 S: (1/2-)22-36/37-45-60-61		
602-046-00-2	heptachlor (ISO); 1,4,5,6,7,8,8-heptachloro-3a,4,7,7a-tetrahydro-4,7-methanoindene	200-962-3	76-44-8	T; R24/25 Carc. Cat. 3; R40 R33 N; R50-53	T; N R: 24/25-33-40-50/53 S: (1/2-)36/37-45-60-61		
602-047-00-8	chlordane (ISO); 1,2,4,5,6,7,8,8-octachloro-3a,4,7,7a-tetrahydro-4,7-methanoindan	200-349-0	57-74-9	Carc. Cat. 3; R40 Xn; R21/22 N; R50-53	Xn; N R: 21/22-40-50/53 S: (2-)36/37-60-61		
602-048-00-3	aldrin (ISO)	206-215-8	309-00-2	T; R24/25-48/24/25 Carc. Cat. 3; R40 N; R50-53	T; N R: 24/25-40-48/24/25-50/53 S: (1/2-)22-36/37-45-60-61		
602-049-00-9	dieldrin (ISO)	200-484-5	60-57-1	T+; R27 T; R25-48/25 Carc. Cat. 3; R40 N; R50-53	T+; N R: 25-27-40-48/25-50/53 S: (1/2-)22-36/37-45-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
602-050-00-4	(1α,4α,4aβ,5β,8β,8aβ)-1,2,3,4,10,10-hexachloro-1,4,4a,5,8,8a-hexahydro-1,4:5,8-dimethanonaphfthalene; isodrin	207-366-2	465-73-6	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)13-28-45-60-61		
602-051-00-X	endrin (ISO); 1,2,3,4,10,10-hexachloro-6,7-epoxy-1,4,4a,5,6,7,8,8a-octahydro-1,4:5,8-dimethanonaphthalene	200-775-7	72-20-8	T+; R28 T; R24 N; R50-53	T+; N R: 24-28-50/53 S: (1/2-)22-36/37-45-60-61		
602-052-00-5	endosulfan (ISO); 1,2,3,4,7,7-hexachloro-8,9,10-trinorborn-2-en-5,6-ylenedimethyl sulphite	204-079-4	115-29-7	T; R24/25 Xi; R36 N; R50-53	T; N R: 24/25-36-50/53 S: (1/2-)28-36/37-45-60-61		
602-053-00-0	isobenzan (ISO); 1,3,4,5,6,7,8,8-octachloro-1,3,3a,4,7,7a-hexahydro-4,7-methanoisobenzofuran	206-045-4	297-78-9	T+; R27/28 N; R50	T+; N R: 27/28-50 S: (1/2-)28-36/37-45-61		
602-054-00-6	3-iodpropene; allyl iodide	209-130-4	556-56-9	R10 C; R34	C R: 10-34 S: (1/2-)7-26-45		
602-055-00-1	bromoethane; ethyl bromide	200-825-8	74-96-4	F; R11 Carc. Cat. 3; R40 Xn; R20/22	F; Xn R: 11-20/22-40 S: (2-)36/37		
602-056-00-7	α,α,α-trifluorotoluene; benzotrifluoride	202-635-0	98-08-8	F; R11 N; R51-53	F; N R: 11-51/53 S: (2-)16-23-61		
602-057-00-2	α-bromotoluene; benzyl bromide	202-847-3	100-39-0	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)39		
602-058-00-8	α,α-dichlorotoluene; benzylidene chloride; benzal chloride	202-709-2	98-87-3	Carc. Cat. 3; R40 T; R23 Xn; R22 Xi; R37/38-41	T R: 22-23-37/38-40-41 S: (1/2-)36/37-38-45		

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602-059-00-3	1-chlorobutane; butyl chloride	203-696-6	109-69-3	F; R11	F R: 11 S: (2-)9-16-29		
602-060-00-9	bromobenzene	203-623-8	108-86-1	R10 Xi; R38 N; R51-53	Xi; N R: 10-38-51/53 S: (2-)61		
602-061-00-4	hexafluoropropene; hexafluoropropylene	204-127-4	116-15-4	Xn; R20 Xi; R37	Xn R: 20-37 S: (2-)41		
602-062-00-X	1,2,3-trichloropropane	202-486-1	96-18-4	Carc. Cat. 2; R45 Repr. Cat. 2; R60 Xn; R20/21/22	T R: 45-60-20/21/22 S: 53-45		E D
602-063-00-5	heptachlor epoxide; 2,3-epoxy-1,4,5,6,7,8,8-heptachloro-3a,4,7,7a-tetrahydro-4,7-methanoindane	213-831-0	1024-57-3	T; R25 Carc. Cat. 3; R40 R33 N; R50-53	T; N R: 25-33-40-50/53 S: (1/2-)36/37-45-60-61		
602-064-00-0	1,3-dichloro-2-propanol	202-491-9	96-23-1	Carc. Cat. 2; R45 T; R25 Xn; R21	T R: 45-21-25 S: 53-45		E
602-065-00-6	hexachlorobenzene	204-273-9	118-74-1	Carc. Cat. 2; R45 T; R48/25 N; R50-53	T; N R: 45-48/25-50/53 S: 53-45-60-61		E
602-066-00-1	tetrachloro-*p*-benzoquinone	204-274-4	118-75-2	Xi; R36/38 N; R50-53	Xi; N R: 36/38-50/53 S: (2-)37-60-61		
602-067-00-7	1,3-dichlorbenzene	208-792-1	541-73-1	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		

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602-068-00-2	ethylene bis(trichloroacetate)	219-732-9	2514-53-6	Xi; R38	Xi R: 38 S: (2-)		
602-069-00-8	dichloroacetylene	—	7572-29-4	E; R2 Carc. Cat. 3; R40 Xn; R48/20	E; Xn R: 2-40-48/20 S: (2-)36/37		
602-070-00-3	3-chloro-4,5,α, α,α-pentafluorotoluene	401-930-3	77227-99-7	R10 Xn; R20/22 N; R50-58	Xn; N R: 10-20/22-50-58 S: (2-)51-60-61		
602-071-00-9	bromobenzylbromotoluene, reaction mass of isomers	402-210-1	99688-47-8	Xn; R48/22 R43 N; R50-53	Xn; N R: 43-48/22-50/53 S: (2-)24-37-41-60-61		
602-072-00-4	dichloro [(dichlorophenyl)methyl]methylbenzene, reaction mass of isomers; (dichlorophenyl)(dichlorotolyl)methane, reaction mass of isomers (IUPAC)	278-404-3	76253-60-6	N; R50-53	N R: 50/53 S: 60-61		
602-073-00-X	1,4-dichlorobut-2-ene	212-121-8	764-41-0	Carc. Cat. 2; R45 T+; R26 T; R24/25 C; R34 N; R50-53	T+; N R: 45-24/25-26-34-50/53 S: 53-45-60-61	Carc. Cat. 2; R45: C ≥ 0,01 % C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	E
602-074-00-5	pentachlorobenzene	210-172-0	608-93-5	F; R11 Xn; R22 N; R50-53	F; Xn; N R: 11-22-50/53 S: (2-)41-46-50-60-61		
602-075-00-0	4,4,5,5-tetrachloro-1,3-dioxolan-2-one	404-060-2	22432-68-4	T+; R26 Xn; R22 C; R34	T+ R: 22-26-34 S: (1/2-)9-26-28-36/37/39-45		

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602-076-00-6	2,3,4-trichlorobut-1-ene	219-397-9	2431-50-7	Carc. Cat. 3; R40 T; R23 Xn; R22 Xi; R36/37/38 N; R50-53	T; N R: 22-23-36/37/38-40-50/53 S: (1/2-)36/37-45-60-61	Carc. Cat. 3; R40: C ≥ 0,1 %	
602-077-00-1	dodecachloropentacyclo[5.2.1.$0^{2,6}$.$0^{3,9}$.$0^{5,8}$]decane; mirex	219-196-6	2385-85-5	Carc. Cat. 3; R40 Repr. Cat. 3; R62-63 R64 Xn; R21/22 N; R50/53	Xn; N R: 21/22-40-50/53-62-63-64 S: (2-)13-36/37-46-60-61		
602-078-00-7	hexachlorocyclopentadiene	201-029-3	77-47-4	T+; R26 T; R24 Xn; R22 C; R34 N; R50-53	T+; N R: 22-24-26-34-50/53 S: (1/2-)25-39-45-53-60-61		
602-079-00-2	2,3-dichloropropene; 2,3-dichloropropylene	201-153-8	78-88-6	F; R11 Muta. Cat. 3; R68 Xn; R20/21/22 Xi; R37/38-41 R52-53	F; Xn R: 11-20/21/22-37/38-41-52/53-68 S: (2-)9-16-23-26-36/37/39-61		
602-080-00-8	alkanes, C_{10-13}, chloro	287-476-5	85535-84-8	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)24-36/37-60-61		
602-081-00-3	2-chloro-4,5-difluorobenzoic acid	405-380-5	—	Xn; R21/22 Xi; R41 R43	Xn R: 21/22-41-43 S: (2-)26-36/37/39		
602-082-00-9	2,2,6,6-tetrakis(bromomethyl)-4-oxaheptane-1,7-diol	408-020-5	109678-33-3	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-41-61		

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602-083-00-4	diphenyl ether, pentabromo derivative pentabromodiphenyl ether	251-084-2	32534-81-9	Xn; R48/21/22 R64 N; R50-53	Xn; N R: 48/21/22-50/53-64 S: (1/2-)36/37-45-60-61		
602-084-00-X	1,1-dichloro-1-fluoroethane	404-080-1	1717-00-6	R52-53 N; R59	N R: 52/53-59 S: 59-61		
602-085-00-5	2-bromopropane	200-855-1	75-26-3	F; R11 Repr. Cat. 1; R60 Xn; R48/20 R66	F; T R: 60-11-48/20-66 S: 53-45		E
602-086-00-0	trifluoroiodomethane; trifluoromethyl iodide	219-014-5	2314-97-8	Muta. Cat. 3; R68	Xn R: 68 S: (2-)36/37		
602-087-00-6	1,2,4-trichlorobenzene	204-428-0	120-82-1	Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-50/53 S: (2-)23-37/39-60-61		
602-088-00-1	2,3-dibromopropan-1-ol; 2,3-dibromo-1-propanol	202-480-9	96-13-9	Carc. Cat. 2; R45 Repr. Cat. 3; R62 T; R24 Xn; R20/22 R52-53	T R: 45-20/22-24-52/53-62 S: 53-45-61		E
602-089-00-7	4-bromo-2-chlorofluorobenzene	405-580-2	60811-21-4	Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-50/53 S: (2-)26-36/37-60-61		
602-090-00-2	1-allyl-3-chloro-4-fluorobenzene	406-630-6	121626-73-1	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)23-37-61		
602-091-00-8	1,3-dichloro-4-fluorobenzene	406-160-1	1435-48-9	Xn; R22-48/20/22 Xi; R38 N; R51-53	Xn; N R: 22-38-48/20/22-51/53 S: (2-)36/37-61		

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602-092-00-3	1-bromo-3,4,5-trifluorobenzene	418-480-9	138526-69-9	R10 Carc. Cat. 3; R40 Xi; R38-41 N; R51-53	Xn; N R: 10-38-40-41-51/53 S: (2-)23-26-36/37/39-61		
602-093-00-9	α, α,α,4-tetrachlorotoluene; *p*-chlorobenzotrichloride	226-009-1	5216-25-1	Carc. Cat. 2; R45 Repr. Cat. 3; R62 T; R48/23 Xn; R21/22 Xi; R37/38	T R: 45-21/22-37/38-48/23-62 S: 53-45		E
602-094-00-4	diphenylether; octabromo derivate	251-087-9	32536-52-0	Repr. Cat. 2; R61 Repr. Cat. 3; R62	T R: 61-62 S: 53-45		
602-096-00-5	malachite green hydrochloride; [1] malachite green oxalate [2]	209-322-8 [1] 219-441-7 [2]	569-64-2 [1] 2437-29-8 [2]	Repr. Cat. 3; R63 Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-63-50/53 S: (2-)26-36/37-39-46-60-61		
602-097-00-0	1-bromo-9-(4,4,5,5,5-pentafluoropentylthio)nonane	422-850-5	148757-89-5	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
603-001-00-X	methanol	200-659-6	67-56-1	F; R11 T; R23/24/25-39/23/24/25	F; T R: 11-23/24/25-39/23/24/25 S: (1/2-)7-16-36/37-45	T; R23/24/25: C ≥ 20 % Xn; R20/21/22: 3 % ≤ C < 20 % T; R39/23/24/25: C ≥ 10 % Xn; R68/20/21/22: 3 % ≤ C < 10 %	
603-002-00-5	ethanol; ethyl alcohol	200-578-6	64-17-5	F; R11	F R: 11 S: (2-)7-16		
603-003-00-0	propan-1-ol; *n*-propanol	200-746-9	71-23-8	F; R11 Xi; R41 R67	F; Xi R: 11-41-67 S: (2-)7-16-24-26-39		

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603-004-00-6	butan-1-ol; *n*-butanol	200-751-6	71-36-3	R10 Xn; R22 Xi; R37/38-41 R67	Xn R: 10-22-37/38-41-67 S: (2-)7/9-13-26-37/39-46		
603-005-00-1	2-methylpropan-2-ol; *tert*-butyl alcohol	200-889-7	75-65-0	F; R11 Xn; R20	F; Xn R: 11-20 S: (2-)9-16		
603-006-00-7	pentanol isomers, with the exception fo those specified elsewhere in this Annex	250-378-8		R10 Xn; R20 Xi; R37 R66	Xn R: 10-20-37-66 S: (2-)46		C
603-007-00-2	2-methylbutan-2-ol; *tert*-pentanol	200-908-9	75-85-4	F; R11 Xn; R20 Xi; R37/38	F; Xn R: 11-20-37/38 S: (2-)46		
603-008-00-8	4-methylpentan-2-ol; methyl isobutyl carbinol	203-551-7	108-11-2	R10 Xi; R37	Xi R: 10-37 S: (2-)24/25	Xi; R37: C ≥ 25 %	
603-009-00-3	cyclohexanol	203-630-6	108-93-0	Xn; R20/22 Xi; R37/38	Xn R: 20/22-37/38 S: (2-)24/25		
603-010-00-9	2-methylcyclohexanol, mixed isomers; [1] *cis*-2-methylcyclohexanol; [2] *trans*-2-methylcyclohexanol [3]	209-512-0 [1] 231-187-9 [2] 231-186-3 [3]	583-59-5 [1] 7443-70-1 [2] 7443-52-9 [3]	Xn; R20	Xn R: 20 S: (2-)24/25		C
603-011-00-4	2-methoxyethanol; ethylene glycol monomethyl ether	203-713-7	109-86-4	R10 Repr. Cat. 2; R60-61 Xn; R20/21/22	T R: 60-61-10-20/21/22 S: 53-45		E
603-012-00-X	2-ethoxyethanol; ethylene glycol monoethyl ether	203-804-1	110-80-5	R10 Repr. Cat. 2; R60-61 Xn; R20/21/22	T R: 60-61-10-20/21/22 S: 53-45		E

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603-013-00-5	2-isopropoxyethanol; ethylene glycol monoisopropyl ether	203-685-6	109-59-1	Xn; R20/21 Xi; R36	Xn R: 20/21-36 S: (2-)24/25		
603-014-00-0	2-butoxyethanol; ethylene glycol monobutyl ether; butyl cellosolve	203-905-0	111-76-2	Xn; R20/21/22 Xi; R36/38	Xn R: 20/21/22-36/38 S: (2-)36/37-46		
603-015-00-6	allyl alcohol	203-470-7	107-18-6	R10 T; R23/24/25 Xi; R36/37/38 N; R50	T; N R: 10-23/24/25-36/37/38-50 S: (1/2-)36/37/39-38-45-61		
603-016-00-1	4-hydroxy-4-methylpentan-2-one; diacetone alcohol	204-626-7	123-42-2	Xi; R36	Xi R: 36 S: (2-)24/25	Xi; R36: C ≥ 10 %	
603-018-00-2	furfuryl alcohol	202-626-1	98-00-0	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)	Xn; R20/21/22: C ≥ 5 %	
603-019-00-8	dimethyl ether	204-065-8	115-10-6	F+; R12	F+ R: 12 S: (2-)9-16-33		
603-020-00-3	ethyl methyl ether	—	540-67-0	F+; R12	F+ R: 12 S: (2-)9-16-33		
603-021-00-9	methyl vinyl ether	203-475-4	107-25-5	F+; R12	F+ R: 12 S: (2-)9-16-33		D
603-022-00-4	diethyl ether; ether	200-467-2	60-29-7	F+; R12 R19 Xn; R22 R66 R67	F+; Xn R: 12-19-22-66-67 S: (2-)9-16-29-33		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-023-00-X	ethylene oxide; oxirane	200-849-9	75-21-8	F+; R12 ⊗ Carc. Cat. 2; R45 Muta. Cat. 2; R46 T; R23 Xi; R36/37/38	F+; T R: 45-46-12-23-36/37/38 S: 53-45		E
603-024-00-5	1,4-dioxane	204-661-8	123-91-1	F; R11-19 Carc. Cat. 3; R40 Xi; R36/37 R66	F; Xn R: 11-19-36/37-40-66 S: (2-)9-16-36/37-46		D
603-025-00-0	tetrahydrofuran	203-726-8	109-99-9	F; R11-19 Xi; R36/37	F; Xi R: 11-19-36/37 S: (2-)16-29-33	Xi; R36/37: C ≥ 25 %	
603-026-00-6	1-chloro-2,3-epoxypropane; epichlorhydrin	203-439-8	106-89-8	R10 Carc. Cat. 2; R45 T; R23/24/25 C; R34 R43	T R: 45-10-23/24/25-34-43 S: 53-45	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,1 % ≤ C < 1 %	E
603-027-00-1	ethanediol; ethylene glycol	203-473-3	107-21-1	Xn; R22	Xn R: 22 S: (2-)		
603-028-00-7	2-chloroethanol; ethylene chlorohydrin	203-459-7	107-07-3	T+; R26/27/28	T+ R: 26/27/28 S: (1/2-)7/9-28-45		
603-029-00-2	bis(2-chloroethyl) ether	203-870-1	111-44-4	R10 ⊗ Carc. Cat. 3; R40 T+; R26/27/28	T+ R: 10-26/27/28-40 S: (1/2-)7/9-27-28-36/37-45		
603-030-00-8	2-aminoethanol; ethanolamine	205-483-3	141-43-5	Xn; R20/21/22 C; R34	C R: 20/21/22-34 S: (1/2-)26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-031-00-3	1,2-dimethoxyethane; ethylene glycol dimethyl ether; EGDME	203-794-9	110-71-4	F; R11 R19 Repr. Cat. 2; R60 Repr. Cat. 2; R61 Xn; R20	F; T R: 60-61-11-19-20 S: 53-45		E
603-032-00-9	ethylene dinitrate; ethylene glycol dinitrate	211-063-0	628-96-6	E; R2 $^{\$}$ T+; R26/27/28 R33	E; T+ R: 2-26/27/28-33 S: (1/2-)33-35-36/37-45		
603-033-00-4	oxydiethylene dinitrate; diethylene glycol dinitrate; digol dinitrate	211-745-8	693-21-0	E; R3 T+; R26/27/28 R33 R52-53	E; T+ R: 3-26/27/28-33-52/53 S: (1/2-)33-35-36/37-45-61		
603-034-00-X	glycerol trinitrate; nitroglycerine	200-240-8	55-63-0	E; R3 T+; R26/27/28 R33 N; R51-53	E; T+; N R: 3-26/27/28-33-51/53 S: (1/2-)33-35-36/37-45-61		
603-035-00-5	pentaerythritol tetranitrate; P.E.T.N.	201-084-3	78-11-5	E; R3	E R: 3 S: (2-)35		
603-036-00-0	mannitol hexanitrate; nitromannite	239-924-6	15825-70-4	E; R3	E R: 3 S: (2-)35		
603-037-00-6	cellulose nitrate; nitrocellulose, containing more than 12,6 % nitrogen	—	—	E; R3 $^{\$}$ R1	E R: 1-3 S: (2-)35		
603-037-01-3	cellulose nitrate; nitrocellulose, containing a maximum of 12,6 % nitrogen	—	—	F; R11 $^{\$}$	F R: 11 S: (2-)16-33-37/39		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-038-00-1	allyl glycidyl ether; allyl 2,3-epoxypropyl ether; prop-2-en-1-yl 2,3-epoxypropyl ether	203-442-4	106-92-3	R10 Carc. Cat. 3; R40 Muta. Cat. 3; R68 Repr. Cat. 3; R62 Xn; R20/22 Xi; R37/38-41 R43 R52-53	Xn R: 10-20/22-37/38-40-41-43-52/53-62-68 S: (2-)24/25-26-36/37/39-61		
603-039-00-7	butyl glycidyl ether; butyl 2,3-epoxypropyl ether	219-376-4	2426-08-6	R10 Carc. Cat. 3; R40 Muta. Cat. 3; R68 Xn; R20/22 Xi; R37 R43 R52-53	Xn R: 10-20/22-37-40-43-52/53-68 S: (2-)24/25-36/37-61		
603-040-00-2	sodium methanolate; sodium methoxide; [1] potassium methanolate; potassium methoxide; [2] lithium methanolate; lithium methoxide [3]	204-699-5 [1] 212-736-1 [2] 212-737-7 [3]	124-41-4 [1] 865-33-8 [2] 865-34-9 [3]	F; R11 C; R34 R14	F; C R: 11-14-34 S: (1/2-)8-16-26-43-45		
603-041-00-8	potassium ethanolate; potassium ethoxide; [1] sodium ethanolate; sodium ethoxide [2]	213-029-0 [1] 205-487-5 [2]	917-58-8 [1] 141-52-6 [2]	F; R11 C; R34 R14	F; C R: 11-14-34 S: (1/2-)8-16-26-43-45		
603-042-00-3	aluminium-tri-isopropoxide	209-090-8	555-31-7	F; R11	F R: 11 S: (2-)8-16		
603-043-00-9	triarimol (ISO); 2,4-dichloro-α-(pyrimidin-5-yl) benzhydryl alcohol	—	26766-27-8	Xn; R22	Xn R: 22 S: (2-)		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-044-00-4	dicofol (ISO); 2,2,2-trichloro-1,1-bis(4-chlorophenyl)ethanol	204-082-0	115-32-2	Xn; R21/22 Xi; R38 R43 N; R50-53	Xn; N R: 21/22-38-43-50/53 S: (2-)36/37-60-61		
603-045-00-X	diisopropyl ether; [1] dipropyl ether [2]	203-560-6 [1] 203-869-6 [2]	108-20-3 [1] 111-43-3 [2]	F; R11 R19 R66 R67	F R: 11-19-66-67 S: (2-)9-16-29-33		C
603-046-00-5	bis (chloromethyl) ether; oxybis(chloromethane)	208-832-8	542-88-1	R10 $ Carc. Cat. 1; R45 T+; R26 T; R24 Xn; R22	T+ R: 45-10-22-24-26 S: 53-45	Carc. Cat. 1; R45: C ≥ 0,001 %	E
603-047-00-0	2-dimethylaminoethanol; *N,N*-dimethylethanolamine	203-542-8	108-01-0	R10 Xn; R20/21/22 C; R34	C R: 10-20/21/22-34 S: (1/2-)25-26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
603-048-00-6	2-diethylaminoethanol; *N,N*-diethylethanolamine	202-845-2	100-37-8	R10 Xn; R20/21/22 C; R34	C R: 10-20/21/22-34 S: (1/2-)25-26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
603-049-00-1	chlorfenethol (ISO); 1,1-bis (4-chlorophenyl) ethanol	201-246-3	80-06-8	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)36-61		
603-050-00-7	1-(2-Butoxypropoxy)propan-2-ol	246-011-6	24083-03-2	Xn; R21/22	Xn R: 21/22 S: (2-)		
603-051-00-2	2-ethylbutan-1-ol	202-621-4	97-95-0	Xn; R21/22	Xn R: 21/22 S: (2-)		
603-052-00-8	3-butoxypropan-2-ol; propylene glycol monobutyl ether	225-878-4	5131-66-8	Xi; R36/38	Xi R: 36/38 S: (2-)		

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603-053-00-3	2-methylpentane-2,4-diol	203-489-0	107-41-5	Xi; R36/38	Xi R: 36/38 S: (2-)	Xi; R36/38: C ≥ 10 %	
603-054-00-9	di-*n*-butyl ether; dibutyl ether	205-575-3	142-96-1	R10 Xi; R36/37/38 R52-53	Xi R: 10-36/37/38-52/53 S: (2-)61	Xi; R36/37/38: C ≥ 10 %	
603-055-00-4	propylene oxide; 1,2-epoxypropane; methyloxirane	200-879-2	75-56-9	F+; R12 Carc. Cat. 2; R45 Muta. Cat. 2; R46 Xn; R20/21/22 Xi; R36/37/38	F+; T R: 45-46-12-20/21/22-36/37/38 S: 53-45		E
603-056-00-X	[(*p*-tolyloxy)methyl]oxirane; [1] [(*m*-tolyloxy)methyl]oxirane; [2] 2,3-epoxypropyl *o*-tolyl ether; [3] [(tolyloxy)methyl]oxirane; cresyl glycidyl ether [4]	218-574-8 [1] 218-575-3 [2] 218-645-3 [3] 247-711-4 [4]	2186-24-5 [1] 2186-25-6 [2] 2210-79-9 [3] 26447-14-3 [4]	Muta. Cat. 3; R68 Xi; R38 R43 N; R51-53	Xn; N R: 38-43-51/53-68 S: (2-)36/37-61		C
603-057-00-5	benzyl alcohol	202-859-9	100-51-6	Xn; R20/22	Xn R: 20/22 S: (2-)26		
603-058-00-0	1,3-propylene oxide	207-964-3	503-30-0	F; R11 Xn; R20/21/22	F; Xn R: 11-20/21/22 S: (2-)9-16-26-29		
603-059-00-6	hexan-1-ol	203-852-3	111-27-3	Xn; R22	Xn R: 22 S: (2-)24/25		

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603-060-00-1	2,2'-bioxirane; 1,2:3,4-diepoxybutane	215-979-1	1464-53-5	Carc. Cat. 2; R45 Muta. Cat. 2; R46 T+; R26 T; R24/25 C; R34	T+ R: 45-46-24/25-26-34 S: 53-45		E
603-061-00-7	tetrahydro-2-furylmethanol; tetrahydrofurfuryl alcohol	202-625-6	97-99-4	Xi; R36	Xi R: 36 S: (2-)39	Xi; R36: C ≥ 10 %	
603-062-00-2	tetrahydrofuran-2,5-diyldimethanol	203-239-0	104-80-3	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)39	Xi; R36/37/38: C ≥ 10 %	
603-063-00-8	2,3-epoxypropan-1-ol; glycidol; oxiranemethanol	209-128-3	556-52-5	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 2; R60 T; R23 Xn; R21/22 Xi; R36/37/38	T R: 45-60-21/22-23-36/37/38-68 S: 53-45		E
603-064-00-3	1-methoxy-2-propanol; monopropylene glycol methyl ether	203-539-1	107-98-2	R10	R: 10 S: (2-)24		
603-065-00-9	resorcinol diglycidyl ether; 1,3-bis(2,3-epoxypropoxy)benzene	202-987-5	101-90-6	Carc. Cat. 3; R40 Muta. Cat. 3; R68 Xn; R21/22 Xi; R36/38 R43 R52-53	Xn R: 21/22-36/38-40-43-52/53-68 S: (2-)23-36/37-61		
603-066-00-4	1,2-epoxy-4-epoxyethylcyclohexane; vinylcyclohexane diepoxide	203-437-7	106-87-6	T; R23/24/25 Xn; R68	T R: 23/24/25-68 S: (1/2-)23-24-45	T; R23/24/25: C ≥ 1 %: Xn; R20/21/22: 0,1 % ≤ C < 1 %	

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603-067-00-X	phenyl glycidyl ether; 2,3-epoxypropyl phenyl ether; 1,2-epoxy-3-phenoxypropane	204-557-2	122-60-1	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R20 Xi; R37/38 R43 R52-53	T R: 45-20-37/38-43-68-52/53 S: 53-45-61		E
603-068-00-5	2,3-epoxypropyl-2-ethylcyclohexyl ether; ethylcyclohexylglycidyl ether	—	130014-35-6	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)26-28-37/39		
603-069-00-0	2,4,6-tris(dimethylaminomethyl)phenol	202-013-9	90-72-2	Xn; R22 Xi; R36/38	Xn R: 22-36/38 S: (2-)26-28		
603-070-00-6	2-amino-2-methylpropanol	204-709-8	124-68-5	Xi; R36/38 R52-53	Xi R: 36/38-52/53 S: (2-)61	Xi; R36/38: C ≥ 10 %	
603-071-00-1	2,2'-iminodiethanol; diethanolamine	203-868-0	111-42-2	Xn; R22-48/22 Xi; R38-41	Xn R: 22-38-41-48/22 S: (2-)26-36/37/39-46		
603-072-00-7	1,4-bis(2,3 epoxypropoxy)butane; butanedioldiglycidyl ether	219-371-7	2425-79-8	Xn; R20/21 Xi; R36/38 R43	Xn R: 20/21-36/38-43 S: (2-)26-28-37/39		
603-073-00-2	bis-[4-(2,3-epoxipropoxi)phenyl]propane	216-823-5	1675-54-3	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)28-37/39	Xi; R36/38: C ≥ 5 %	
603-074-00-8	reaction product: bisphenol-A-(epichlorhydrin); epoxy resin (number average molecular weight ≤ 700)	500-033-5	25068-38-6	Xi; R36/38 R43 N; R51-53	Xi; N R: 36/38-43-51/53 S: (2-)28-37/39-61	Xi; R36/38: C ≥ 5 %	
603-075-00-3	chlormethyl methyl ether; chlorodimethyl ether	203-480-1	107-30-2	F; R11 Carc. Cat. 1; R45 Xn; R20/21/22	F; T R: 45-11-20/21/22 S: 53-45		E

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603-076-00-9	but-2-yne-1,4-diol; 2-butyne-1,4-diol	203-788-6	110-65-6	C; R34 T; R23/25 Xn; R21-48/22 R43	C; T R: 21-23/25-34-43-48/22 S: (1/2-)25-26-36/37/39-45-46	C; R34: C ≥ 50 % Xi; R36/38: 25 % ≤ C < 50 %	D
603-077-00-4	1-dimethylaminopropan-2-ol; dimepranol (INN)	203-556-4	108-16-7	R10 Xn; R22 C; R34	C R: 10-22-34 S: (1/2-)23-26-36-45		
603-078-00-X	prop-2-yn-1-ol; propargyl alcohol	203-471-2	107-19-7	R10 T; R23/24/25 C; R34 N; R51-53	T; N R: 10-23/24/25-34-51/53 S: (1/2-)26-28-36-45-61		
603-079-00-5	2,2'-(methylimino)diethanol; N-methyldiethanolamine	203-312-7	105-59-9	Xi; R36	Xi R: 36 S: (2-)24		
603-080-00-0	2-methylaminoethanol; N-methylethanolamine; N-methyl-2-ethanolamine; N-methyl-2-amino ethanol; 2-(methylamino)ethanol	203-710-0	109-83-1	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
603-081-00-6	2,2'-thiodiethanol; thiodiglycol	203-874-3	111-48-8	Xi; R36	Xi R: 36 S: (2-)		
603-082-00-1	1-aminopropan-2-ol; isopropanolamine	201-162-7	78-96-6	C; R34	C R: 34 S: (1/2-)23-26-36-45		
603-083-00-7	1,1'-iminodipropan-2-ol; di-isopropanolamine	203-820-9	110-97-4	Xi; R36	Xi R: 36 S: (2-)26		
603-084-00-2	styrene oxide; (epoxyethyl)benzene; phenyloxirane	202-476-7	96-09-3	Carc. Cat. 2; R45 Xn; R21 Xi; R36	T R: 45-21-36 S: 53-45		E

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603-085-00-8	bronopol (INN); 2-bromo-2-nitropropane-1,3-diol	200-143-0	52-51-7	Xn; R21/22 Xi; R37/38-41 N; R50	Xn; N R: 21/22-37/38-41-50 S: (2-)26-37/39-61		
603-086-00-3	ethirimol (ISO); 5-butyl-2-ethylamino-6-methylpyrimidin-4-ol	245-949-3	23947-60-6	Xn; R21	Xn R: 21 S: (2-)36/37		
603-087-00-9	2-ethylhexane-1,3-diol; octylene glycol; ethoexadiol	202-377-9	94-96-2	Xi; R41	Xi R: 41 S: (2-)25-26-39-46		
603-088-00-4	2-(octylthio)ethanol; 2-hydroxyethyl octyl sulphide	222-598-4	3547-33-9	Xi; R41	Xi R: 41 S: (2-)26		
603-089-00-X	7,7-dimethyl-3-oxa-6-azaoctan-1-ol	400-390-6	—	C; R35 Xn; R22	C R: 22-35 S: (1/2-)26-28-36/37/39-45		
603-090-00-5	2-(2-bromoethoxy)anisole	402-010-4	4463-59-6	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
603-091-00-0	*exo*-1-methyl-4-(1-methylethyl)-7-oxabicyclo[2.2.1]heptan-2-ol	402-470-6	87172-89-2	Xn; R22 Xi; R41	Xn R: 22-41 S: (2-)26-39		
603-092-00-6	2-methyl-4-phenylpentanol	402-770-7	92585-24-5	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
603-093-00-1	cinmethylin (ISO) *exo*-(±)-1-methyl-2-(2-methylbenzyloxy)-4-isopropyl-7-oxabicyclo(2.2.1)heptane	402-410-9	87818-31-3	Xn; R20 N; R51-53	Xn; N R: 20-51/53 S: (2-)23-61		
603-094-00-7	1,3-bis(2,3-epoxypropoxy)-2,2-dimethylpropane	241-536-7	17557-23-2	Xi; R38 R43	Xi R: 38-43 S: (2-)24-37		

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603-095-00-2	2-(propyloxy)ethanol; EGPE	220-548-6	2807-30-9	Xn; R21 Xi; R36	Xn R: 21-36 S: (2-)26-36/37-46		
603-096-00-8	2-(2-butoxyethoxy)ethanol; diethylene glycol monobutyl ether	203-961-6	112-34-5	Xi; R36	Xi R: 36 S: (2-)24-26		
603-097-00-3	1,1',1"-nitrilotripropan-2-ol; triisopropanolamine	204-528-4	122-20-3	Xi; R36 R52-53	Xi R: 36-52/53 S: (2-)26-61		
603-098-00-9	2-phenoxyethanol	204-589-7	122-99-6	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)26		
603-099-00-4	3-(*N*-methyl-*N*-(4-methylamino-3-nitrophenyl)amino)propane-1,2-diol hydrochloride	403-440-5	93633-79-5	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
603-100-00-8	1,2-dimethoxypropane	404-630-0	7778-85-0	F; R11-19	F R: 11-19 S: (2-)9-16-24/25-33		
603-101-00-3	tetrahydro-2-isobutyl-4-methylpyran-4-ol, mixed isomers (*cis* and *trans*)	405-040-6	—	Xi; R36	Xi R: 36 S: (2-)25-26		
603-102-00-9	1,2-epoxybutane	203-438-2	106-88-7	F; R11 Carc. Cat. 3; R40 Xn; R20/21/22 Xi; R36/37/38 R52-53	F; Xn R: 11-20/21/22-36/37/38-40-52/53 S: (2-)9-16-29-36/37-61		
603-103-00-4	oxirane, mono[(C_{12-14}-alkyloxy)methyl] derivs.	271-846-8	68609-97-2	Xi; R38 R43	Xi R: 38-43 S: (2-)24-37		

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603-104-00-X	fenarimol (ISO); 2,4'-dichloro-α-(pyrimidin-5-yl)benzhydryl alcohol	262-095-7	60168-88-9	Repr. Cat. 3; R62-63 R64 N; R51-53	Xn; N R: 51/53-62-63-64 S: (2-)36/37-61		
603-105-00-5	furan	203-727-3	110-00-9	F+; R12 R19 Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R20/22-48/22 Xi; R38 R52-53	F+; T R: 45-12-19-20/22-38-48/22-68-52/53 S: 53-45-61		E
603-106-00-0	2-methoxypropanol	216-455-5	1589-47-5	R10 Repr. Cat. 2; R61 Xi; R37/38-41	T R: 61-10-37/38-41 S: 53-45		
603-107-00-6	2-(2-methoxyethoxy)ethanol; diethylene glycol monomethyl ether	203-906-6	111-77-3	Repr. Cat. 3; R63	Xn R: 63 S: (2-)36/37		
603-108-00-1	2-methylpropan-1-ol; iso-butanol	201-148-0	78-83-1	R10 Xi; R37/38-41 R67	Xi R: 10-37/38-41-67 S: (2-)7/9-13-26-37/39-46		
603-117-00-0	propan-2-ol; isopropyl alcohol; isopropanol	200-661-7	67-63-0	F; R11 Xi; R36 R67	F; Xi R: 11-36-67 S: (2-)7-16-24/25-26		
603-118-00-6	6-dimethylaminohexan-1-ol	404-680-3	1862-07-3	Xn; R22 C; R34 R52-53	C R: 22-34-52/53 S: (1/2-)26-36/37/39-45-61		
603-119-00-1	1,1'-(1,3-phenylenedioxy)bis(3-(2-(prop-2-enyl)phenoxy)propan-2-ol)	405-840-5	—	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-120-00-7	2-methyl-5-phenylpentanol	405-890-8	25634-93-9	Xi; R36/38	Xi R: 36/38 S: (2-)26-37		
603-121-00-2	4-[4-(1,3-dihydroxyprop-2-yl)phenylamino]-1,8-dihydroxy-5-nitroanthraquinone	406-057-1	114565-66-1	Carc. Cat. 3; R40 R43 R53	Xn R: 40-43-53 S: (2-)36/37-61		
603-122-00-8	sodium 2-ethylhexanolate	406-150-7	38411-13-1	F; R11 C; R34 R52-53	F; C R: 11-34-52/53 S: (1/2-)7-26-36/37/39-45-61		
603-123-00-3	4-methyl-8-methylenetricyclo[3.3.1.1^{3,7}]decan-2-ol	406-330-5	122760-84-3	Xi; R38 R43 N; R51-53	Xi; N R: 38-43-51/53 S: (2-)24-37-61		
603-124-00-9	1,4-bis[2-(vinyloxy)ethoxy]benzene	406-900-3	84563-49-5	N; R50-53	N R: 50/53 S: 60-61		
603-125-00-4	2-(2,4-dichlorophenyl)-1-(1*H*-1,2,4-triazol-1-yl)pent-4-en-2-ol	407-850-5	89544-40-1	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)26-39-61		
603-126-00-X	2-((4-methyl-2-nitrophenyl)amino)ethanol	408-090-7	100418-33-5	Xn; R22 R43 R52-53	Xn R: 22-43-52/53 S: (2-)36/37-61		
603-127-00-5	butan-2-ol; [1] (*S*)-butan-2-ol; [2] (*R*)-butan-2-ol; [3] (±)-butan-2-ol [4]	201-158-5 [1] 224-168-1 [2] 238-967-8 [3] 240-029-8 [4]	78-92-2 [1] 4221-99-2 [2] 14898-79-4 [3] 15892-23-6 [4]	R10 Xi; R36/37 R67	Xi R: 10-36/37-67 S: (2-)7/9-13-24/25-26-46		C
603-128-00-0	2-(phenylmethoxy)naphthalene	405-490-3	613-62-7	R53	R: 53 S: 61		
603-129-00-6	1-*tert*-butoxypropan-2-ol	406-180-0	57018-52-7	R10 Xi; R41	Xi R: 10-41 S: (2-)26-39		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-130-00-1	reaction mass of isomers of: α-((dimethyl)biphenyl)-ω-hydroxypoly(oxyethylene)	406-325-8	—	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)39-61		
603-131-00-7	reaction mass of: 1-deoxy-1-[methyl-(1-oxododecyl)amino]-D-glucitol; 1-deoxy-1-[methyl-(1-oxotetradecyl)amino]-D-glucitol (3:1)	407-290-1	—	Xi; R41	Xi R: 41 S: (2-)26-39		
603-132-00-2	2-hydroxymethyl-9-methyl-6-(1-methylethyl)-1,4-dioxaspiro [4.5]decane	408-200-3	63187-91-7	Xi; R38-41 R52-53	Xi R: 38-41-52/53 S: (2-)26-37/39-61		
603-133-00-8	reaction mass of: 3-[(4-amino-2-chloro-5-nitrophenyl) amino]-propane-1,2-diol; 3,3'-(2-chloro-5-nitro-1,4-phenylenediimino)bis(propan-1,2-diol)	408-240-1	—	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-36-61		
603-134-00-3	reaction mass of substituted dodecyl and/or tetradecyl, diphenyl ethers. The substance is produced by the Friedel Crafts reaction. The catalyst is removed from the reaction product. Diphenyl ether is substituted by C_1-C_{10} alkyl groups. The alkyl groups are bonded randomly between C_1 and C_6. Linear C_{12} and C_{14}, 50/50 used.	410-450-3	—	R53	R: 53 S: 61		
603-135-00-9	bis[[2,2',2"-nitrilotris-[ethanolato]]-1-N,O]-bis[2-(2-methoxyethoxy)ethoxy]-titanium	410-500-4	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
603-136-00-4	3-((4-(bis(2-hydroxyethyl)amino)-2-nitrophenyl)amino)-1-propanol	410-910-3	104226-19-9	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
603-137-00-X	reaction mass of: 1-deoxy-1-[methyl-(1-oxohexadecyl)amino]-D-glucitol; 1-deoxy-1-[methyl-(1-oxooctadecyl)amino]-D-glucitol	411-130-6	—	Xi; R41	Xi R: 41 S: (2-)26-39		
603-138-00-5	3-(2,2-dimethyl-3-hydroxypropyl)toluene; (alt.): 2,2-dimethyl-3-(3-methylphenyl)propanol	403-140-4	103694-68-4	R52-53	R: 52/53 S: 61		
603-139-00-0	bis(2-methoxyethyl) ether	203-924-4	111-96-6	R10 R19 Repr. Cat. 2; R60-61	T R: 60-61-10-19 S: 53-45		

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603-140-00-6	2,2' -oxybisethanol; diethylene glycol	203-872-2	111-46-6	Xn; R22	Xn R: 22 S: (2-)46		
603-141-00-1	reaction mass of: dodecyloxy-1-methyl-1-[oxy-poly-(2-hydroxymethylethanoxy)]pentadecane; dodecyloxy-1-methyl-1-[oxy-poly-(2-hydroxymethylethanoxy)]heptadecane	413-780-6	—	R52-53	R: 52/53 S: 61		
603-142-00-7	2-(2-(2-hydroxyethoxy)ethyl)-2-aza-bicyclo[2.2.1]heptane	407-360-1	116230-20-7	Xn; R21/22-48/20 Xi; R38-41	Xn R: 21/22-38-41-48/20 S: (2-)26-36/37/39		
603-143-00-2	*R*-2,3-epoxy-1-propanol	404-660-4	57044-25-4	E; R2 Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 2; R60 T; R23 Xn; R21/22 C; R34	E; T R: 45-60-2-21/22-23-34-68 S: 53-45		E
603-144-00-8	reaction mass of: 2,6,9-trimethyl-2,5,9-cyclododecatrien-1-ol; 6,9-dimethyl-2-methylen-5,9-cyclododecadien-1-ol	413-530-6	111850-00-1	N; R51-53	N R: 51/53 S: 61		
603-145-00-3	2-isopropyl-2-(1-methylbutyl)-1,3-dimethoxypropane	406-970-5	129228-11-1	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)36/37-61		
603-146-00-9	2-[(2-[2-(dimethylamino)ethoxy]ethyl)methylamino]ethanol	406-080-7	83016-70-0	Xn; R22 C; R34 R52-53	C R: 22-34-52/53 S: (1/2-)23-26-36/37/39-45-61		
603-147-00-4	(-)-*trans*-4-(4'-fluorophenyl)-3-hydroxymethyl-*N*-methylpiperidine	406-030-4	105812-81-5	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)22-24-26-37/39-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-148-00-X	1,4-bis[(vinyloxy)methyl]cyclohexane	413-370-7	17351-75-6	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
603-149-00-5	reaction mass of diastereoisomers of 1-(1-hydroxyethyl)-4-(1-methylethyl)cyclohexane	407-640-3	63767-86-2	Xi; R36/38 N; R51-53	Xi; N R: 36/38-51/53 S: (2-)26-37-61		
603-150-00-0	(±) *trans*-3,3-dimethyl-5-(2,2,3-trimethyl-cyclopent-3-en-1-yl)-pent-4-en-2-ol	411-580-3	107898-54-4	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)24/25-37-60-61		
603-151-00-6	(±)-2-(2,4-dichlorophenyl)-3-(1*H*-1,2,4-triazol-1-yl)propan-1-ol	413-570-4	—	R52-53	R: 52/53 S: 61		
603-152-00-1	2-(4-*tert*-butylphenyl)ethanol	410-020-5	5406-86-0	Repr. Cat. 3; R62 Xn; R48/22 Xi; R41 N; R51-53	Xn; N R: 41-48/22-62-51/53 S: (2-)26-36/37/39-61		
603-153-00-7	3-((2-nitro-4-(trifluoromethyl)phenyl)amino)propane-1,2-diol	410-010-0	104333-00-8	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
603-154-00-2	1-[(2-*tert*-butyl)cyclohexyloxy]-2-butanol	412-300-2	139504-68-0	N; R51-53	N R: 51/53 S: 61		
603-155-00-8	Reaction products of 2-(4,6-bis(2,4-dimethylphenyl)-1,3,5-triazin-2-yl)-5-hydroxyphenol with ((C$_{10-16}$, rich in C$_{12-13}$ alkyloxy)methyl)oxyrane	410-560-1	—	N; R50-53	N R: 50/53 S: 60-61		
603-156-00-3	2-(2,4-dichlorophenyl)-2-(2-propenyl)oxirane	411-210-0	89544-48-9	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
603-157-00-9	6,9-bis(hexadecyloxymethyl)-4,7-dioxanonane-1,2,9-triol	411-450-6	143747-72-2	R53	R: 53 S: 61		
603-158-00-4	reaction mass of 4 diastereoisomers of 2,7-dimethyl-10-(1-methylethyl)-1-oxaspiro[4.5]deca-3,6-diene	412-460-3	—	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		

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603-159-00-X	2-cyclododecylpropan-1-ol	411-410-8	118562-73-5	N; R50-53	N R: 50/53 S: 60-61		
603-160-00-5	1,2-diethoxypropane	412-180-1	10221-57-5	F; R11 R19	F R: 11-19 S: (2-)9-16-24-33		
603-161-00-0	1,3-diethoxypropane	413-140-6	3459-83-4	R10	R: 10 S: (2-)9-24		
603-162-00-6	α[2-[[[(2-hydroxyethyl)methylamino]acetyl]amino]propyl]-ω-(nonylphenoxy)poly[oxo(methyl-1,2-ethanediyl)]	413-420-8	144736-29-8	C; R34 R43 N; R51-53	C; N R: 34-43-51/53 S: (1/2-)26-28-36/37/39-45-61		
603-163-00-1	2-phenyl-1,3-propanediol	411-810-2	1570-95-2	Xi; R41	Xi R: 41 S: (2-)26-39		
603-164-00-7	2-butyl-4-chloro-4,5-dihydro-5-hydroxymethyl-1-[2'-(2-triphenylmethyl-1,2,3,4-2*H*-tetrazol-5-yl)-1,1'-biphenyl-4-methyl]-1*H*-imidazole	412-420-5	133909-99-6	R53	R: 53 S: 61		
603-165-00-2	reaction mass of: 4-allyl-2,6-bis(2,3-epoxypropyl)phenol; 4-allyl-6-[3-[6-[3-[6-[3-(4-allyl-2,6-bis(2,3-epoxypropyl)phenoxy)-2-hydroxypropyl]-4-allyl-2-(2,3-epoxypropyl)phenoxy]-2-hydroxypropyl]-4-allyl-2-(2,3-epoxypropyl)phenoxy]-2-hydroxypropyl]-2-(2,3-epoxypropyl)phenol; 4-allyl-6-[3-(4-allyl-2,6-bis(2,3-epoxypropyl)phenoxy)-2-hydroxypropyl]-2-(2,3-epoxypropyl)phenol; 4-allyl-6-[3-[6-[3-(4-allyl-2,6-bis(2,3-epoxypropyl)phenoxy)-2-hydroxypropyl]-4-allyl-2-(2,3-epoxypropyl)phenoxy]-2-hydroxypropyl]-2-(2,3-epoxypropyl)phenol	417-470-1	—	Muta. Cat. 3; R68 R43	Xn R: 43-68 S: (2-)36/37		
603-166-00-8	*R*-1-chloro-2,3-epoxypropane	424-280-2	51594-55-9	R10 Carc. Cat. 2; R45 T; R23/24/25 C; R34 R43	T R: 45-10-23/24/25-34-43 S: 53-45		E
603-167-00-3	3,3',5,5'-tetra-*tert*-butylbiphenyl-2,2'-diol	407-920-5	6390-69-8	R53	R: 53 S: 61		

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603-168-00-9	3-(2-ethylhexyloxy)propane-1,2-diol	408-080-2	70445-33-9	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
603-169-00-4	(±)-*trans*-4-(4-fluorophenyl)-3-hydroxymethyl-*N*-methylpiperidine	415-550-0	109887-53-8	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)22-26-39-61		
603-170-00-X	reaction mass of: 2-methyl-1-(6-methylbicyclo[2.2.1]hept-5-en-2-yl)pent-1-en-3-ol; 2-methyl-1-(1-methylbicyclo[2.2.1]hept-5-en-2-yl)-pent-1-en-3-ol; 2-methyl-1-(5-methylbicyclo[2.2.1]hept-5-en-2-yl)pent-1-en-3-ol	415-990-3	67739-11-1	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
603-171-00-5	5-thiazolylmethanol	414-780-9	38585-74-9	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
603-172-00-0	mono-2-[2-(4-dibenzo[*b*,*f*][1,4]thiazepin-11-yl)piperazinium-1-yl]ethoxy)ethanol *trans*-butenedioate	415-180-1	773058-82-5	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)22-26-39-61		
603-173-00-6	4,4-dimethyl-3,5,8-trioxabicyclo[5.1.0]octane	421-750-9	57280-22-5	Xi; R36 R43	Xi R: 36-43 S: (2-)26-36/37		
603-174-00-1	4-cyclohexyl-2-methyl-2-butanol	420-630-3	83926-73-2	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
603-175-00-7	2-(2-hexyloxyethoxy)ethanol; DEGHE; diethylene glycol monohexyl ether; 3,6-dioxa-1-dodecanol; hexyl carbitol; 3,6-dioxadodecan-1-ol	203-988-3	112-59-4	Xn; R21 Xi; R41	Xn R: 21-41 S: (2-)26-36/37/39-46		
603-176-00-2	1,2-bis(2-methoxyethoxy)ethane; TEGDME; triethylene glycol dimethyl ether; triglyme	203-977-3	112-49-2	R19 Repr. Cat. 2; R61 Repr. Cat. 3; R62	T R: 61-19-62 S: 53-45		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-177-00-8	1-ethoxypropan-2-ol; 2PG1EE; 1-ethoxy-2-propanol; propylene glycol monoethyl ether; [1] 2-ethoxy-1-methylethyl acetate; 2PG1EEA [2]	216-374-5 [1] 259-370-9 [2]	1569-02-4 [1] 54839-24-6 [2]	R10 R67	R: 10-67 S: (2-)24		
603-178-00-3	2-hexyloxyethanol; ethylene glycol monohexyl ether; n-hexylglycol	203-951-1	112-25-4	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-36/37/39-45		
603-179-00-9	ergocalciferol (ISO); Vitamin D2	200-014-9	50-14-6	T+; R26 T; R24/25-48/25	T+ R: 24/25-26-48/25 S: (1/2-)28-36/37-45		
603-180-00-4	colecalciferol; Vitamin D3	200-673-2	67-97-0	T+; R26 T; R24/25-48/25	T+ R: 24/25-26-48/25 S: (1/2-)28-36/37-45		
603-181-00-X	*tert*-butyl methyl ether; MTBE; 2-methoxy-2-methylpropane	216-653-1	1634-04-4	F; R11 Xi; R38	F; Xi R: 11-38 S: (2-)9-16-24		
603-183-00-0	2-[2-(2-butoxyethoxy)ethoxy]ethanol; TEGBE; triethylene glycol monobutyl ether; butoxytriethylene glycol	205-592-6	143-22-6	Xi; R41	Xi R: 41 S: (2-)26-39-46	Xi; R41: C ≥ 30 % Xi; R36: 20 % ≤ C < 30 %	
603-184-00-6	2-(hydroxymethyl)-2-[[2-hydroxy-3-(isooctadecyloxy)propoxy]methyl]-1,3-propanediol	416-380-1	146925-83-9	N; R50-53	N R: 50/53 S: 60-61		
603-185-00-1	2,4-dichloro-3-ethyl-6-nitrophenol	420-740-1	99817-36-4	T; R25 Xi; R41 R43 N; R50-53	T; N R: 25-41-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
603-186-00-7	trans-(5*RS*,6*SR*)-6-amino-2,2-dimethyl-1,3-dioxepan-5-ol	419-050-3	79944-37-9	R43	Xi R: 43 S: (2-)22-24/25-26-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
603-187-00-2	2-((4,6-bis(4-(2-(1-methylpyridinium-4-yl)vinyl)phenylamino)-1,3,5-triazin-2-yl)(2-hydroxyethyl)amino)ethanol dichloride	419-360-9	163661-77-6	N; R50-53	N R: 50/53 S: 60-61		
603-189-00-3	reaction mass of complexes of: titanium, 2,2'-oxydiethanol, ammonium lactate, nitrilotris(2-propanol) and ethylene glycol	405-250-8	—	N; R51-53	N R: 51/53 S: 61		
603-191-00-4	2-(4,6-bis(2,4-dimethylphenyl)-1,3,5-triazin-2-yl)-5-(3-((2-ethylhexyl)oxy)-2-hydroxypropoxy)phenol	419-740-4	137658-79-8	R53	R: 53 S: 61		
603-195-00-6	2-[4-(4-methoxyphenyl)-6-phenyl-1,3,5-triazin-2-yl]-phenol	430-810-3	154825-62-4	R52-53	R: 52/53 S: 61		
603-196-00-1	2-(7-ethyl-1*H*-indol-3-yl)ethanol	431-020-1	41340-36-7	Xn; 22-48/22 N; R51-53	Xn; N R: 22-48/22-51/53 S: (2-)36/37/39-61		
603-197-00-7	tebuconazole (ISO); 1-(4-chlorophenyl)-4,4-dimethyl-3-(1,2,4-triazol-1-ylmethyl) pentan-3-ol	403-640-2	107534-96-3	Repr. Cat. 3; R63 Xn; R22 N; R51-53	Xn; N R: 22-51/53-63 S: (2-)22-36/37-61		
603-199-00-8	etoxazol (ISO); (*RS*)-5-tert-butyl-2-[2-(2,6-difluorophenyl)-4,5-dihydro-1,3-oxazol-4-yl]phenetole	—	153233-91-1	N; R50-53	N R: 50/53 S: 60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
604-001-00-2	phenol; carbolic acid; monohydroxybenzene; phenylalcohol	203-632-7	108-95-2	Muta. Cat. 3; R68 T; R23/24/25 Xn; R48/20/21/22 C; R34	T; R: 23/24/25-34-48/20/21/22-68 S: (1/2-)24/25-26-28-36/37/39-45	T; R23/24/25: C ≥ 10 % Xn; R20/21/22: 3 % ≤ C < 10 % C; R34: C ≥ 3 % Xi; R36/38: 1 % ≤ C < 3 %	
604-002-00-8	pentachlorophenol	201-778-6	87-86-5	Carc. Cat. 3; R40 T+; R26 T; R24/25 Xi; R36/37/38 N; R50-53	T+; N R: 24/25-26-36/37/38-40-50/53 S: (1/2-)22-36/37-45-52-60-61		

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604-003-00-3	sodium pentachlorophenolate; [1] potassium pentachlorophenolate [2]	205-025-2 [1] 231-911-3 [2]	131-52-2 [1] 7778-73-6 [2]	Carc. Cat. 3; R40 T+; R26 T; R24/25 Xi; R36/37/38 N; R50-53	T+; N R: 24/25-26-36/37/38-40-50/53 S: (1/2-)22-28-36/37-45-52-60-61		
604-004-00-9	*m*-cresol; [1] *o*-cresol; [2] *p*-cresol; [3] mix-cresol [4]	203-577-9 [1] 202-423-8 [2] 203-398-6 [3] 215-293-2 [4]	108-39-4 [1] 95-48-7 [2] 106-44-5 [3] 1319-77-3 [4]	T; R24/25 C; R34	T R: 24/25-34 S: (1/2-)36/37/39-45	T; R24/25: C ≥ 5 % Xn; R21/22: 1 % ≤ C < 5 % C; R34: C ≥ 5 % Xi; R36/38: 1 % ≤ C < 5 %	C
604-005-00-4	1,4-dihydroxybenzene; hydroquinone; quinol	204-617-8	123-31-9	Carc. Cat. 3; R40 Muta. Cat. 3; R68 Xn; R22 Xi; R41 R43 N; R50	Xn; N R: 22-40-41-43-50-68 S: (2-)26-36/37/39-61		
604-006-00-X	3,4-xylenol; [1] 2,5-xylenol; [2] 2,4-xylenol; [3] 2,3-xylenol; [4] 2,6-xylenol; [5] xylenol; [6] 2,4(or 2,5)-xylenol [7]	202-439-5 [1] 202-461-5 [2] 203-321-6 [3] 208-395-3 [4] 209-400-1 [5] 215-089-3 [6] 276-245-4 [7]	95-65-8 [1] 95-87-4 [2] 105-67-9 [3] 526-75-0 [4] 576-26-1 [5] 1300-71-6 [6] 71975-58-1 [7]	T; R24/25 C; R34 N; R51-53	T; N R: 24/25-34-51/53 S: (1/2-)26-36/37/39-45-61		C
604-007-00-5	2-naphthol	205-182-7	135-19-3	Xn; R20/22 N; R50	Xn; N R: 20/22-50 S: (2-)24/25-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
604-008-00-0	2-chlorophenol; [1] 4-chlorophenol; [2] 3-chlorophenol; [3] chlorophenol [4]	202-433-2 [1] 203-402-6 [2] 203-582-6 [3] 246-691-4 [4]	95-57-8 [1] 106-48-9 [2] 108-43-0 [3] 25167-80-0 [4]	Xn; R20/21/22 N; R51-53	Xn; N R: 20/21/22-51/53 S: (2-)28-61		C
604-009-00-6	pyrogallol; 1,2,3-trihydroxybenzene	201-762-9	87-66-1	Muta. Cat. 3; R68 Xn; R20/21/22 R52-53	Xn R: 20/21/22-68-52/53 S: (2-)36/37-61	Xn; R20/21/22: C ≥ 10 %	
604-010-00-1	resorcinol; 1,3-benzenediol	203-585-2	108-46-3	Xn; R22 Xi; R36/38 N; R50	Xn; N R: 22-36/38-50 S: (2-)26-61	Xn; R22: C ≥ 10 %	
604-011-00-7	2,4-dichlorophenol	204-429-6	120-83-2	T; R24 Xn; R22 C; R34 N; R51-53	T; N R: 22-24-34-51/53 S: (1/2-)26-36/37/39-45-61		
604-012-00-2	4-chloro-*o*-cresol; 4-chloro-2-methyl phenol	216-381-3	1570-64-5	T; R23 C; R35 N; R50	T; C; N R: 23-35-50 S: (1/2-)26-36/37/39-45-61	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
604-013-00-8	2,3,4,6-tetrachlorophenol	200-402-8	58-90-2	T; R25 Xi; R36/38 N; R50-53	T; N R: 25-36/38-50/53 S: (1/2-)26-28-37-45-60-61	T; R25: C ≥ 5 % Xn; R22: 0,5 % ≤ C < 5 % Xi; R36/38: C ≥ 5 %	
604-014-00-3	chlorocresol; 4-chloro-*m*-cresol; 4-chloro-3-methylphenol	200-431-6	59-50-7	Xn; R21/22 Xi; R41 R43 N; R50	Xn; N R: 21/22-41-43-50 S: (2-)26-36/37/39-61	Xn; R21/22: C ≥ 10 %	
604-015-00-9	2,2'-methylenebis-(3,4,6-trichlorophenol); hexachlorophene	200-733-8	70-30-4	T; R24/25 N; R50-53	T; N R: 24/25-50/53 S: (1/2-)20-37-45-60-61	T; R24/25: C ≥ 2 % Xn; R21/22: 0,2 % ≤ C < 2 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
604-016-00-4	1,2-dihydroxybenzene; pyrocatechol	204-427-5	120-80-9	Xn; R21/22 Xi; R36/38	Xn R: 21/22-36/38 S: (2-)22-26-37		
604-017-00-X	2,4,5-trichlorophenol	202-467-8	95-95-4	Xn; R22 Xi; R36/38 N; R50-53	Xn; N R: 22-36/38-50/53 S: (2-)26-28-60-61	Xn; R22: C ≥ 20 % Xi; R36/38: C ≥ 5 %	
604-018-00-5	2,4,6-trichlorophenol	201-795-9	88-06-2	Carc. Cat. 3; R40 Xn; R22 Xi; R36/38 N; R50-53	Xn; N R: 22-36/38-40-50/53 S: (2-)36/37-60-61		
604-019-00-0	dichlorophen (ISO)	202-567-1	97-23-4	Xn; R22 Xi; R36 N; R50-53	Xn; N R: 22-36-50/53 S: (2-)26-60-61		
604-020-00-6	2-phenylphenol (ISO); biphenyl-2-ol; 2-hydroxybiphenyl;	201-993-5	90-43-7	Xi; R36/37/38 N; R50	Xi; N R: 36/37/38-50 S: (2-)22-61		
604-021-00-1	sodium 2-biphenylate; 2-phenylphenol, sodium salt	205-055-6	132-27-4	Xn; R22 Xi; R37/38-41 N; R50	Xn; N R: 22-37/38-41-50 S: (2-)22-26-61		
604-022-00-7	2,2-dimethyl-1,3-benzodioxol-4-ol	400-900-7	22961-82-6	Xi; R41	Xi R: 41 S: (2-)24-26-39		
604-023-00-2	2,4-dichloro-3-ethylphenol	401-060-4	—	C; R34 N; R50-53	C; N R: 34-50/53 S: (1/2-)26-36/39-45-60-61		
604-024-00-8	4,4-isobutylethylidenediphenol	401-720-1	6807-17-6	Repr. Cat. 2; R60 Xi; R36 N; R50-53	T; N R: 60-36-50/53 S: 53-45-60-61		
604-025-00-3	2,5-bis(1,1-dimethylbutyl)hydroquinone	400-220-0	—	N; R51-53	N R: 51/53 S: 61		

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604-026-00-9	2,2-spirobi(6-hydroxy-4,4,7-trimethylchromane)	400-270-3	—	N; R51-53	N R: 51/53 S: 61		
604-027-00-4	2-methyl-5-(1,1,3,3-tetramethylbutyl)hydroquinone	400-530-6	—	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24/25-26-37-61		
604-028-00-X	4-amino-3-fluorophenol	402-230-0	399-95-1	Carc. Cat. 2; R45 Xn; R22 R43 N; R51-53	T; N R: 45-22-43-51/53 S: 53-45-61		E
604-029-00-5	1-naphtol	201-969-4	90-15-3	Xn; R21/22 Xi; R37/38-41	Xn R: 21/22-37/38-41 S: (2-)22-26-37/39		
604-030-00-0	bisphenol A; 4,4'-isopropylidenediphenol	201-245-8	80-05-7	Repr. Cat. 3; R62 Xi; R37-41 R43	Xn R: 37-41-43-62 S: (2-)26-36/37-39-46		
604-031-00-6	guaiacol	201-964-7	90-05-1	Xn; R22 Xi; R36/38	Xn R: 22-36/38 S: (2-)26		
604-032-00-1	thymol	201-944-8	89-83-8	Xn; R22 C; R34 N; R51-53	C; N R: 22-34-51/53 S: (1/2-)26-28-36/37/39-45-61		
604-033-00-7	isobutyl but-3-enoate	401-170-2	24342-03-8	R10	R: 10 S: (2-)		
604-034-00-2	4,4'-thiodi-*o*-cresol	403-330-7	24197-34-0	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
604-035-00-8	4-nonylphenol, reaction products with formaldehyde and dodecane-1-thiol	404-160-6	—	R43 R53	X R: 43-53 S: (2-)24-37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
604-036-00-3	4,4'-oxybis(ethylenethio)diphenol	404-590-4	90884-29-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
604-037-00-9	3,5-xylenol; 3,5-dimethylphenol	203-606-5	108-68-9	T; R24/25 C; R34	T R: 24/25-34 S: (1/2-)26-28-36/37/39-45		
604-038-00-4	4-chloro-3,5-dimethylphenol; [1] chloroxylenol [2]	201-793-8 [1] 215-316-6 [2]	88-04-0 [1] 1321-23-9 [2]	Xn; R22 Xi; R36/38 R43	Xn R: 22-36/38-43 S: (2-)24-37		
604-039-00-X	ethyl 2-[4-[(6-chlorobenzoxazol-2-yl)oxy]phenoxy]propionate; fenoxaprop-ethyl	266-362-9	66441-23-4	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
604-040-00-5	fomesafen (ISO); 5-[2-chloro-4-(trifluoromethyl)phenoxy]-*N*-(methylsulphonyl)-2-nitrobenzamide	276-439-9	72178-02-0	Xn; R22	Xn R: 22 S: (2-)		
604-041-00-0	acifluorfen (ISO); 5-[2-chloro-4-(trifluoromethyl)phenoxy]-2-nitrobenzoic acid [1] sodium 5-[2-chloro-4-(trifluoromethyl) phenoxy]-2-nitrobenzoate; acifluorfen-sodium [2]	256-634-5 [1] 263-560-7 [2]	50594-66-6 [1] 62476-59-9 [2]	Xn; R22 Xi; R38-41 N; R50-53	Xn; N R: 22-38-41-50/53 S: (2-)24-39-60-61		
604-042-00-6	4-nitrosophenol	203-251-6	104-91-6	Muta. Cat. 3; R68 Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53-68 S: (2-)26-36/37/39-47-49-61		
604-043-00-1	monobenzone; 4-hydroxyphenyl benzyl ether; hydroquinone monobenzyl ether	203-083-3	103-16-2	Xi; R36 R43	Xi R: 36-43 S: (2-)24/25-26-37		
604-044-00-7	mequinol; 4-methoxyphenol; hydroquinone monomethyl ether	205-769-8	150-76-5	Xn; R22 Xi; R36 R43	Xn R: 22-36-43 S: (2-)24/25-26-37/39-46		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
604-045-00-2	2,3,5-trimethylhydroquinone	211-838-3	700-13-0	Xn; R20 Xi; R37/38-41 R43 N; R50-53	Xn; N R: 20-37/38-41-43-50/53 S: (2-)24-26-37/39-60-61		
604-046-00-8	4-(4-isopropoxyphenylsulfonyl)phenol	405-520-5	95235-30-6	N; R51-53	N R: 51/53 S: 61		
604-047-00-3	4-(4-tolyloxy)biphenyl	405-730-7	51601-57-1	Xn; R48/22 R53	Xn R: 48/22-53 S: (2-)22-36-61		
604-048-00-9	4,4',4"-(ethan-1,1,1-triyl)triphenol	405-800-7	27955-94-8	N; R51-53	N R: 51/53 S: 61		
604-049-00-4	4-4'-methylenebis(oxyethylenethio)diphenol	407-480-4	93589-69-6	N; R51-53	N R: 51/53 S: 61		
604-051-00-5	3,5-bis((3,5-di-*tert*-butyl-4-hydroxy)benzyl)-2,4,6-trimethyl-phenol	401-110-5	87113-78-8	R52-53	R: 52/53 S: 61		
604-052-00-0	2,2'-methylenebis(6-(2*H*-benzotriazol-2-yl)-4-(1,1,3,3-tetra-methylbutyl)phenol)	403-800-1	103597-45-1	R53	R: 53 S: 61		
604-053-00-6	2-methyl-4-(1,1-dimethylethyl)-6-(1-methyl-pentadecyl)-phe-nol	410-760-9	157661-93-3	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
604-054-00-1	reaction mass of: 2-methoxy-4-(tetrahydro-4-methylene-2*H*-pyran-2-yl)-phenol; 4-(3,6-dihydro-4-methyl-2*H*-pyran-2-yl)-2-methoxyphenol	412-020-0	—	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
604-055-00-7	2,2'-((3,3',5,5'-tetramethyl-(1,1'-biphenyl)-4,4'-diyl)-bis(oxy-methylene))-bis-oxirane	413-900-7	85954-11-6	Muta. Cat. 3; R68	Xn R: 68 S: (2-)22-36-37		
604-056-00-2	2-(2-hydroxy-3,5-dinitroanilino)ethanol	412-520-9	99610-72-7	F; R11 Repr. Cat. 3; R62 Xn; R22	F; Xn R: 11-22-62 S: (2-)22-33-36/37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
604-057-00-8	reaction mass of: isomers of 2-(2*H*-benzotriazol-2-yl)-4-methyl-(*n*)-dodecylphenol; isomers of 2-(2*H*-benzotriazol-2-yl)-4-methyl-(*n*)-tetracosyl-phenol; isomers of 2-(2*H*-benzotriazol-2-yl)-4-methyl-5,6-didodecyl-phenol. *n* = 5 or 6	401-680-5	—	N; R51-53	N R: 51/53 S: 61		
604-058-00-3	1,2-bis(3-methylphenoxy)ethane	402-730-9	54914-85-1	N; R50-53	N R: 50/53 S: 60-61		
604-059-00-9	2-*n*-hexadecylhydroquinone	406-400-5	—	Xn; R48/22 Xi; R38 R43 R53	Xn R: 38-43-48/22-53 S: (2-)22-36/37-61		
604-060-00-4	9,9-bis(4-hydroxyphenyl)fluorene	406-950-6	3236-71-3	Xi; R36-38 N; R50-53	Xi; N R: 36/38-50/53 S: (2-)26-37-60-61		
604-061-00-X	reaction mass of: 2-chloro-5-*sec*-tetradecylhydroquinones where *sec*-tetradecyl = 1-methyltridecyl; 1-ethyldodecyl; 1-propylundecyl; 1-butyldecyl; 1-pentylnonyl; 1-hexyloctyl	407-740-7	—	Xi; R38 R43 R52-53	Xi R: 38-43-52/53 S: (2-)24-37-61		
604-062-00-5	2,4-dimethyl-6-(1-methyl-pentadecyl)phenol	411-220-5	—	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
604-063-00-0	5,6-dihydroxyindole	412-130-9	3131-52-0	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)22-26-36/37/39-61		
604-064-00-6	2-(4,6-diphenyl-1,3,5-triazin-2-yl)-5-((hexyl)oxy)-phenol	411-380-6	147315-50-2	R53	R: 53 S: 61		
604-065-00-1	4,4',4"-(1-methylpropan-1-yl-3-ylidene)tris(2-cyclohexyl-5-methylphenol)	407-460-5	111850-25-0	N; R51-53	N R: 51/53 S: 61		

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604-066-00-7	reaction mass of: phenol, 6-(1,1-dimethylethyl)-4-tetrapropyl-2-[(2-hydroxy-5-tetra-propylphenyl)methyl (C_{41}-compound) and methane, 2,2'-bis[6-(1,1-dimethyl-ethyl)-1-hydroxy-4-tetrapropyl-phenyl)]- (C_{45}-compound); 2,6-bis(1,1-dimethylethyl)-4-tetra-propyl-phenol and 2-(1,1-dimethylethyl)-4-tetrapropyl-phenol; 2,6-bis[(6-(1,1-dimethylethyl)-1-hydroxy-4-tetrapropylphenyl) methyl]-4-(tetrapropyl)phenol and 2-[(6-(1,1-dimethylethyl)-1-hydroxy-4-tetrapropylphenylmethyl]-6-[1-hydroxy-4-tetra-propylphenyl)methyl]-4-(tetrapropyl)phenol	414-550-8	—	N; R50-53	N R: 50/53 S: 60-61		

604-067-00-2	reaction mass of: 2,2'-[[(2-hydroxyethyl)imino]bis(methylene) bis[4-dodecylphenol];
	formaldehyde, oligomer with 4-dodecyl phenol and 2-aminoethanol(n = 2);
	formaldehyde, oligomer with 4-dodecyl phenol and 2-aminoethanol(n = 3, 4 and higher)
		414-520-4	—	Xi; R38-41
				N; R50-53
					Xi; N
					R: 38-41-50/53
					S: (2-)26-37/39-60-61		

604-068-00-8	(±)-4-[2-[[3-(4-hydroxyphenyl)-1-methylpropyl]amino]-1-hydroxyethyl]phenol hydrochloride	415-170-5	90274-24-1	Xn; R20/22 R43	Xn R: 20/22-43 S: (2-)24-26-37		
604-069-00-3	2-(1-methylpropyl)-4-*tert*-butylphenol	421-740-4	51390-14-8	C; R34 N; R51-53	C; N R: 34-51/53 S: (1/2-)26-36/37/39-45-61		
604-070-00-9	triclosan; 2,4,4'-trichloro-2'-hydroxy-diphenyl-ether; 5-chloro-2-(2,4-dichlorophenoxy)phenol	222-182-2	3380-34-5	Xi; R36/38 N; R50-53	Xi; N R: 36/38-50/53 S: 26-39-46-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
605-001-00-5	formaldehyde … %	200-001-8	50-00-0	Carc. Cat. 3; R40 T; R23/24/25 C; R34 R43	T R: 23/24/25-34-40-43 S: (1/2-)26-36/37/39-45-51	T; R23/24/25: C ≥25 % Xn; R20/21/22: 5 % ≤ C < 25 % C; R34: C ≥25 % Xi; R36/37/38: 5 % ≤ C < 25 % R43: C ≥ 0,2 %	B D
605-002-00-0	1,3,5-trioxan; trioxymethylene	203-812-5	110-88-3	F; R11 Repr. Cat. 3; R63 Xi; R37	F; Xn R: 11-37-63 S: (2-)36/37-46		

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605-003-00-6	acetaldehyde; ethanal	200-836-8	75-07-0	F+; R12 Carc. Cat. 3; R40 Xi; R36/37	F+; Xn R: 12-36/37-40 S: (2-)16-33-36/37		
605-004-00-1	2,4,6-trimethyl-1,3,5-trioxan; paraldehyde	204-639-8	123-63-7	F; R11 $	F R: 11 S: (2-)9-16-29-33		
605-005-00-7	2,4,6,8-tetramethyl-1,3,5,7-tetraoxacyclooctane; metaldehyde	203-600-2	108-62-3	R10 $ Xn; R22	Xn R: 10-22 S: (2-)13-25-46		
605-006-00-2	butyraldehyde	204-646-6	123-72-8	F; R11	F R: 11 S: (2-)9-29-33		
605-007-00-8	1,1-dimethoxyethane; dimethyl acetal	208-589-8	534-15-6	F; R11	F R: 11 S: (2-)9-16-33		
605-008-00-3	acrylaldehyde; acrolein; prop-2-enal	203-453-4	107-02-8	F; R11 T+; R26 T; R24/25 C; R34 N; R50	F; T+; N R: 11-24/25-26-34-50 S: 23-26-28-36/37/39-45-61		D
605-009-00-9	crotonaldehyde; 2-butenal; [1] (*E*)-2-butenal; (*E*)-crotonaldehyde [2]	224-030-0 [1] 204-647-1 [2]	4170-30-3 [1] 123-73-9 [2]	F; R11 Muta. Cat. 3; R68 T+; R26 T; R24/25 Xn; R48/22 Xi; R37/38-41 N; R50	F; T+; N R: 11-24/25-26-37/38-41-48/22-50-68 S: (1/2-)26-28-36/37/39-45-61		
605-010-00-4	2-furaldehyde	202-627-7	98-01-1	Carc. Cat. 3; R40 T; R23/25 Xn; R21 Xi; R36/37	T R: 21-23/25-36/37-40 S: (1/2-)26-36/37/39-45	T; R23/25: C ≥ 5 % Xn; R20/22: 1 % ≤ C < 5 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
605-011-00-X	2-chlorobenzaldehyde; *o*-chlorobenzaldehyde	201-956-3	89-98-5	C; R34	C R: 34 S: (1/2-)26-45		
605-012-00-5	benzaldehyde	202-860-4	100-52-7	Xn; R22	Xn R: 22 S: (2-)24		
605-013-00-0	chloralose (INN); (*R*)-1,2-O-(2,2,2-trichloroethylidene)-α-D-glucofuranose; glucochloralose; anhydroglucochloral	240-016-7	15879-93-3	Xn; R20/22	Xn R: 20/22 S: (2-)16-24/25-28		
605-014-00-6	chloral hydrate; 2,2,2-trichloroethane-1,1-diol	206-117-5	302-17-0	T; R25 Xi; R36/38	T R: 25-36/38 S: (1/2-)25-45		
605-015-00-1	1,1-diethoxyethane; acetal	203-310-6	105-57-7	F; R11 Xi; R36/38	F; Xi R: 11-36/38 S: (2-)9-16-33	Xi; R36/38: C ≥ 10 %	
605-016-00-7	glyoxal … %; ethandial … %	203-474-9	107-22-2	Muta. Cat. 3; R68 Xn; R20 Xi; R36/38 R43	Xn R: 20-36/38-43-68 S: (2-)36/37	Xn; R20: C ≥ 10 % Xi; R36/38: C ≥ 10 %	B
605-017-00-2	1,3-dioxolane	211-463-5	646-06-0	F; R11	F R: 11 S: (2-)16		
605-018-00-8	propanal; propionaldehyde	204-623-0	123-38-6	F; R11 Xi; R36/37/38	F; Xi R: 11-36/37/38 S: (2-)9-16-29		
605-019-00-3	citral	226-394-6	5392-40-5	Xi; R38 R43	Xi R: 38-43 S: (2-)24/25-37		
605-020-00-9	safrole; 5-allyl-1,3-benzodioxole	202-345-4	94-59-7	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R22	T R: 45-22-68 S: 53-45		E

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
605-021-00-4	formaldehyde, reaction products with butylphenol	294-145-9	91673-30-2	R43	Xi R: 43 S: (2-)24-37		
605-022-00-X	glutaral; glutaraldehyde; 1,5-pentanedial	203-856-5	111-30-8	T; R23/25 C; R34 R42/43 N; R50	T; N R: 23/25-34-42/43-50 S: (1/2-)26-36/37/39-45-61	T; R25: C ≥ 50 % Xn; R22: 2 % ≤ C < 50 % T; R23: C ≥ 25 % Xn; R20: 2 % ≤ C < 25 % C; R34: C ≥ 10 % Xi; R37/38-41: 2 % ≤ C < 10 % Xi; R36/37/38: 0,5 % ≤ C < 2 % R43: C ≥ 0,5 %	
605-025-00-6	chloroacetaldehyde	203-472-8	107-20-0	Carc. Cat. 3; R40 T+; R26 T; R24/25 C; R34 N; R50	T+; N R: 24/25-26-34-40-50 S: (1/2-)26-28-36/37/39-45-61		
605-026-00-1	2,5,7,7-tetramethyloctanal	405-690-0	114119-97-0	Xi; R38 R43 N; R51-53	Xi; N R: 38-43-51/53 S: (2-)24-37-61		
605-027-00-7	reaction mass of: 3a,4,5,6,7,7a-hexahydro-4,7-methano-1*H*-indene-6-carboxaldehyde; 3a,4,5,6,7,7a-hexahydro-4,7-methano-1*H*-indene-5-carboxaldehyde	410-480-7	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
605-028-00-2	β-methyl-3-(1-methylethyl)-benzenepropanal	412-050-4	125109-85-5	N; R51-53	N R: 51/53 S: 61		
605-029-00-8	2-cyclohexylpropanal	412-270-0	2109-22-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
605-030-00-3	1-(*p*-methoxyphenyl)acetaldehyde oxime	411-510-1	3353-51-3	R43	Xi R: 43 S: (2-)24-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
605-031-00-9	reaction mass of: 2,2-dimethoxyethanal [(this component is considered to be anhydrous in terms of identity, structure and composition. However, 2,2-dimethoxyethanal will exist in a hydrated form. 60 % anhydrous is equivalent to 70.4 % hydrate; water(Including free water and water in hydrated 2,2-dimethoxyethanal)]	421-890-0	—	R43	Xi R: 43 S: (2-)24-37		
606-001-00-8	acetone; propan-2-one; propanone	200-662-2	67-64-1	F; R11 Xi; R36 R66 R67	F; Xi R: 11-36-66-67 S: (2-)9-16-26		
606-002-00-3	butanone; ethyl methyl ketone	201-159-0	78-93-3	F; R11 Xi; R36 R66 R67	F; Xi R: 11-36-66-67 S: (2-)9-16		
606-003-00-9	heptan-3-one; butyl ethyl ketone	203-388-1	106-35-4	R10 Xn; R20 Xi; R36	Xn R: 10-20-36 S: (2-)24		
606-004-00-4	4-methylpentan-2-one; isobutyl methyl ketone	203-550-1	108-10-1	F; R11 Xn; R20 Xi; R36/37 R66	F; Xn R: 11-20-36/37-66 S: (2-)9-16-29		
606-005-00-X	2,6-dimethylheptan-4-one; di-isobutyl ketone	203-620-1	108-83-8	R10 Xi; R37	Xi R: 10-37 S: (2-)24	Xi; R37: C ≥ 10 %	
606-006-00-5	pentan-3-one; diethyl ketone	202-490-3	96-22-0	F; R11 Xi; R37 R66 R67	F; Xi R: 11-37-66-67 S: (2-)9-16-25-33		
606-007-00-0	3-methylbutan-2-one; methyl isopropyl ketone	209-264-3	563-80-4	F; R11	F R: 11 S: (2-)9-16-33		
606-009-00-1	4-methylpent-3-en-2-one; mesityl oxide	205-502-5	141-79-7	R10 Xn; R20/21/22	Xn R: 10-20/21/22 S: (2-)25	Xn; R20/21/22: C ≥ 5 %	

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606-010-00-7	cyclohexanone	203-631-1	108-94-1	R10 Xn; R20	Xn R: 10-20 S: (2-)25		
606-011-00-2	2-methylcyclohexanone	209-513-6	583-60-8	R10 Xn; R20	Xn R: 10-20 S: (2-)25		
606-012-00-8	3,5,5-trimethylcyclohex-2-enone; isophorone	201-126-0	78-59-1	Carc. Cat. 3; R40 Xn; R21/22 Xi; R36/37	Xn R: 21/22-36/37-40 S: (2-)13-23-36/37/39-46	Xi; R36/37: C ≥ 10 %	
606-013-00-3	*p*-benzoquinone; quinone	203-405-2	106-51-4	T; R23/25 Xi; R36/37/38 N; R50	T; N R: 23/25-36/37/38-50 S: (1/2-)26-28-45-61		
606-014-00-9	chlorophacinone (ISO); 2-(2-(4-chlorophenyl)phenylacetyl)indan-1,3-dione	223-003-0	3691-35-8	T+; R27/28 T; R23-48/24/25 N; R50-53	T+; N R: 23-27/28-48/24/25-50/53 S: (1/2-)36/37-45-60-61		
606-016-00-X	pindone (ISO); 2-pivaloylindan-1,3-dione	201-462-8	83-26-1	T; R25-48/25 N; R50-53	T; N R: 25-48/25-50/53 S: (1/2-)37-45-60-61		
606-017-00-5	diketene; diketen	211-617-1	674-82-8	R10 Xn; R20	Xn R: 10-20 S: (2-)3		D
606-018-00-0	dichlone (ISO); 2,3-dichloro-1,4-naphthoquinone	204-210-5	117-80-6	Xn; R22 Xi; R36/38 N; R50-53	Xn; N R: 22-36/38-50/53 S: (2-)26-60-61		
606-019-00-6	chlordecone (ISO); perchloropentacyclo[5,3,0,$0^{2,6}$,$0^{3,9}$,$0^{4,8}$]decan-5-one; decachloropentacyclo[5,2,1,$0^{2,6}$,$0^{3,9}$,$0^{5,8}$]decan-4-one	205-601-3	143-50-0	Carc. Cat. 3; R40 T; R24/25 N; R50-53	T; N R: 24/25-40-50/53 S: (1/2-)22-36/37-45-60-61		
606-020-00-1	5-methylheptan-3-one	208-793-7	541-85-5	R10 Xi; R36/37	Xi R: 10-36/37 S: (2-)23	Xi; R36/37: C ≥ 10 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
606-021-00-7	N-methyl-2-pyrrolidone	212-828-1	872-50-4	Xi; R36/38	Xi R: 36/38 S: (2-)41	Xi; R36/38: C ≥ 10 %	
606-022-00-2	1-phenyl-3-pyrazolidone	202-155-1	92-43-3	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
606-023-00-8	4-methoxy-4-methylpentan-2-one	203-512-4	107-70-0	R10 Xn; R20	Xn R: 10-20 S: (2-)23-24/25		
606-024-00-3	heptan-2-one; methyl amyl ketone	203-767-1	110-43-0	R10 Xn; R20/22	Xn R: 10-20/22 S: (2-)24/25		
606-025-00-9	cyclopentanone	204-435-9	120-92-3	R10 Xi; R36/38	Xi R: 10-36/38 S: (2-)23		
606-026-00-4	5-methylhexan-2-one; isoamyl methyl ketone	203-737-8	110-12-3	R10 Xn; R20	Xn R: 10-20 S: (2-)23-24/25		
606-027-00-X	heptan-4-one; di-*n*-propyl ketone	204-608-9	123-19-3	R10 Xn; R20	Xn R: 10-20 S: (2-)24/25		
606-028-00-5	2,4-dimethylpentan-3-one; di-isopropyl ketone	209-294-7	565-80-0	F; R11 Xn; R20	F; Xn R: 11-20 S: (2-)9-16-24/25		
606-029-00-0	pentane-2,4-dione; acetylacetone	204-634-0	123-54-6	R10 Xn; R22	Xn R: 10-22 S: (2-)21-23-24/25		
606-030-00-6	hexan-2-one; methyl butyl ketone; butyl methyl ketone; methyl-*n*-butyl ketone	209-731-1	591-78-6	R10 Repr. Cat. 3; R62 T; R48/23 R67	T R: 10-48/23-62-67 S: (1/2-)36/37-45		

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606-031-00-1	3-propanolide; 1,3-propiolactone	200-340-1	57-57-8	Carc. Cat. 2; R45 T+; R26 Xi; R36/38	T+ R: 45-26-36/38 S: 53-45		E
606-032-00-7	hexachloroacetone	204-129-5	116-16-5	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)24/25-61		
606-033-00-2	2-(3,4-dichlorophenyl)-4-methyl-1,2,4-oxadiazolidinedione; methazole	243-761-6	20354-26-1	Xn; R21/22 Xi; R36/38 N; R51-53	Xn; N R: 21/22-36/38-51/53 S: (2-)36/37-61		
606-034-00-8	metribuzin (ISO); 4-amino-6-*tert*-butyl-3-methylthio-1,2,4-triazin-5(4*H*)-one; 4-amino-4,5-dihydro-6-(1,1-dimethylethyl)-3-methylthio-1,2,4-triazin-5-one	244-209-7	21087-64-9	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
606-035-00-3	chloridazon (ISO); 5-amino-4-chloro-2-phenylpyridazine-3-(2*H*)-one; pyrazon	216-920-2	1698-60-8	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
606-036-00-9	quinomethionate; chinomethionat (ISO); 6-methyl-1,3-dithiolo(4,5-*b*)quinoxalin-2-one	219-455-3	2439-01-2	Repr. Cat. 3; R62 Xn; R20/21/22-48/22 Xi; R36 R43 N; R50-53	Xn; N R: 20/21/22-36-43-48/22-50/53-62 S: (2-)24-37-60-61		
606-037-00-4	triadimefon (ISO); 1-(4-chlorophenoxy)-3,3-dimethyl-1-(1,2,4-triazol-1-yl)butanone	256-103-8	43121-43-3	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)24-37-61		
606-038-00-X	diphacinone (ISO); 2-diphenylacetylindan-1,3-dione	201-434-5	82-66-6	T+; R28 T; R48/23/24/25	T+ R: 28-48/23/24/25 S: (1/2-)36/37-45		
606-039-00-5	5(or 6)-*tert*-butyl-2'-chloro-6'-ethylamino-3',7'-dimethylspiro(isobenzofuran-1(1*H*),9'-xanthene)-3-one	400-680-2	—	Xn; R20 N; R50-53	Xn; N R: 20-50/53 S: (2-)60-61		

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606-040-00-0	(*N*-benzyl-*N*-ethyl)amino-3-hydroxyacetophenone hydrochloride	401-840-4	55845-90-4	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
606-041-00-6	2-methyl-1-(4-methylthiophenyl)-2-morpholinopropan-1-one	400-600-6	71868-10-5	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-61		
606-042-00-1	acetophenone	202-708-7	98-86-2	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)26		
606-043-00-7	2,4-di-*tert*-butylcyclohexanone	405-340-7	13019-04-0	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
606-044-00-2	2,4,6-trimethylbenzophenone	403-150-9	954-16-5	Xn; R22 Xi; R36 N; R50-53	Xn; N R: 22-36-50/53 S: (2-)26-60-61		
606-045-00-8	oxadiazon (ISO); 3-[2,4-dichloro-5-(1-methylethoxy)phenyl]-5-(1,1-dimethylethyl)-1,3,4-oxadiazol-2(3*H*)-one	243-215-7	19666-30-9	N; R50-53	N R: 50/53 S: 60-61		
606-046-00-3	reaction mass of *cis*- and *trans*-cyclohexadec-8-en-1-one	401-700-2	3100-36-5	N; R50-53	N R: 50/53 S: 60-61		
606-047-00-9	2-benzyl-2-dimethylamino-4-morpholinobutyrophenone	404-360-3	119313-12-1	N; R50-53	N R: 50/53 S: 60-61		
606-048-00-4	2'-anilino-3'-methyl-6'-dipentylaminospiro(isobenzofuran-1(1*H*),9'-xanthen)-3-one	406-480-1	—	R53	R: 53 S: 61		
606-049-00-X	4-(*trans*-4-propylcyclohexyl)acetophenone	406-700-6	78531-61-0	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
606-050-00-5	6-anilino-1-benzoyl-4-(4-*tert*-pentylphenoxy)naphto[1,2,3-de]quinoline-2,7-(3*H*)-dione	412-480-2	72453-58-8	N; R51-53	N R: 51/53 S: 61		

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606-051-00-0	4-pentylcyclohexanone	406-670-4	61203-83-6	N; R51-53	N R: 51/53 S: 61		
606-052-00-6	4-(*N*,*N*-dibutylamino)-2-hydroxy-2'-carboxybenzophenone	410-410-5	54574-82-2	R52-53	R: 52/53 S: 61		
606-053-00-1	flurtamone (ISO); (*RS*)-5-methylamino-2-phenyl-4-(α, α,α-trifluoro-*m*-tolyl) furan-3(2*H*)-one	—	96525-23-4	N; R50-53	N R: 50/53 S: 60-61		
606-054-00-7	isoxaflutole (ISO); 5-cyclopropyl-1,2-oxazol-4-yl α, α,α-trifluoro-2-mesyl-*p*-tolyl ketone	—	141112-29-0	Repr. Cat. 3; R63 N; R50-53	Xn; N R: 50/53-63 S: (2-)36/37-60-61		
606-055-00-2	1-(2,3-dihydro-1,3,3,6-tetramethyl-1-(1-methylethyl)-1*H*-inden-5-yl)ethanone	411-180-9	92836-10-7	Xn; R22-48/22 N; R51-53	Xn; N R: 22-48/22-51/53 S: (2-)24-36-61		
606-056-00-8	4-chloro-3',4'-dimethoxybenzophenone	404-610-1	116412-83-0	N; R50-53	N R: 50/53 S: 60-61		
606-057-00-3	4-propylcyclohexanone	406-810-4	40649-36-3	Xi; R38 R52-53	Xi R: 38-52/53 S: (2-)25-37-61		
606-058-00-9	4'-fluoro-2,2-dimethoxyacetophenone	407-500-1	21983-80-2	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
606-059-00-4	2,4-difluoro-α-(1*H*-1,2,4-triazol-1-yl)acetophenone hydrochloride	412-390-3	86386-75-6	Xn; R22 Xi; R41 R43	Xn R: 22-41-43 S: (2-)22-26-36/37/39		
606-060-00-X	reaction mass of: *trans*-2,4-dimethyl-2-(5,6,7,8-tetrahydro-5,5,8,8-tetramethyl-naphthalene-2-yl)-1,3-dioxolane; *cis*-2,4-dimethyl-2-(5,6,7,8-tetrahydro-5,5,8,8-tetramethyl-naphthalene-2-yl)-1,3-dioxolane	412-950-7	—	N; R50-53	N R: 50/53 S: 60-61		
606-061-00-5	(3-chlorophenyl)-(4-methoxy-3-nitrophenyl)methanone	423-290-4	66938-41-8	Muta. Cat. 3; R68 N; R50-53	Xn; N R: 68-50/53 S: (2-)22-36/37-60-61		

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606-062-00-0	tetrahydrothiopyran-3-carboxaldehyde	407-330-8	61571-06-0	Repr. Cat. 2; R61 Xi; R41 R52-53	T R: 61-41-52/53 S: 53-45-61		
606-063-00-6	(*E*)-3-(2-chlorophenyl)-2-(4-fluorophenyl)propenal	410-980-5	112704-51-5	Xi; R36 R43	Xi R: 36-43 S: (2-)24-26-37		
606-064-00-1	pregn-5-ene-3,20-dione bis(ethylene ketal)	407-450-0	7093-55-2	R53	R: 53 S: 61		
606-065-00-7	1-(4-morpholinophenyl)butan-1-one	413-790-0	—	N; R51-53	N R: 51/53 S: 61		
606-066-00-2	(*E*)-5[(4-chlorophenyl)methylene]-2,2-dimethylcyclopentanone	410-440-9	164058-20-2	N; R51-53	N R: 51/53 S: 61		
606-067-00-8	reaction mass of: 1-(2,3,6,7,8,9-hexahydro-1,1-dimethyl-1*H*-benz(g)inden-4-yl)ethanone; 1-(2,3,5,6,7,8-hexahydro-1,1-dimethyl-1*H*-benz(f)inden-4-yl) ethanone; 1-(2,3,6,7,8,9-hexahydro-1,1-dimethyl-1*H*-benz(g)inden-5-yl) ethanone; 1-(2,3,6,7,8,9-hexahydro-3,3-dimethyl-1*H*-benz(g)inden-5-yl) ethanone	414-870-8	96792-67-5	N; R50-53	N R: 50/53 S: 60-61		
606-068-00-3	2,7,11-trimethyl-13-(2,6,6-trimethylcyclohex-1-en-1-yl)tridecahexaen-2,4,6,8,10,12-al	415-770-7	1638-05-7	Xn; R48/22 R43 R52-53	Xn R: 43-48/22-52/53 S: (2-)22-36/37-61		
606-069-00-9	spiro[1,3-dioxolane-2,5'-(4',4',8',8'-tetramethyl-hexahydro-3',9'-methanonaphthalene)]	415-460-1	154171-76-3	N; R51-53	N R: 51/53 S: 24-61		
606-070-00-4	butroxydim (ISO); 5-(3-butyryl-2,4,6-trimethylphenyl)-2-[1-(ethoxyimino)propyl]-3-hydroxycyclohex-2-en-1-one	414-790-3	138164-12-2	Repr. Cat. 3; R62-63 Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-62-63-50/53 S: (2-)22-36/37-60-61		

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606-071-00-X	17-spiro(5,5-dimethyl-1,3-dioxan-2-yl)androsta-1,4-diene-3-one	421-050-3	13258-43-0	N; R50-53	N R: 50/53 S: 22-60-61		
606-072-00-5	3-acetyl-1-phenyl-pyrrolidine-2,4-dione	421-600-2	719-86-8	Xn; R48/22 N; R51-53	Xn; N R: 48/22-51/53 S: (2-)22-36/37-61		
606-073-00-0	4,4'-bis(dimethylamino)benzophenone; Michler's ketone	202-027-5	90-94-8	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xi; R41	T R: 45-41-68 S: 53-45		
606-075-00-1	1-benzyl-5-ethoxyimidazolidine-2,4-dione	417-340-4	65855-02-9	Xn; R22	Xn R: 22 S: (2-)22		
606-076-00-7	1-((2-quinolinyl-carbonyl)oxy)-2,5-pyrrolidinedione	418-630-3	136465-99-1	Xi; R41 R43	Xi R: 41-43 S: (2-)24-26-37/39		
606-077-00-2	(3*S*,4*S*)-3-hexyl-4-[(*R*)-2-hydroxytridecyl]-2-oxetanone	418-650-2	104872-06-2	N; R50-53	N R: 50/53 S: 60-61		
606-078-00-8	1-octylazepin-2-one	420-040-6	59227-88-2	C; R34 R43 N; R51-53	C; N R: 34-43-51/53 S: (1/2-)26-36/37/39-45-61		
606-079-00-3	2-*n*-butyl-benzo[*d*]isothiazol-3-one	420-590-7	4299-07-4	C; R34 R43 N; R50-53	C; N R: 34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
606-080-00-9	Reaction product of: 3-hydroxy-5,7-di-*tert*-butylbenzofuran-2-one with o-xylene	417-100-9	—	R53	R: 53 S: 61		
606-081-00-4	(3β, 5α, 6β)-3-(acetyloxy)-5-bromo-6-hydroxy-androstan-17-one	419-790-7	4229-69-0	R43 R52-53	Xi R: 43-52/53 S: (2-)22-36/37-61		

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606-082-00-X	reaction mass of: butan-2-one oxime; syn-O,O'-di(butan-2-one oxime)diethoxysilane	406-930-7		T; R48/25 R43 R52-53	T R: 43-48/25-52/53 S: (1/2-)25-36/37-45-61		
606-083-00-5	2-chloro-5-*sec*-hexadecylhydroquinone	407-750-1	137193-60-3	Xi; R36/38 R43 R52-53	Xi R: 36/38-43-52/53 S: (2-)24-26-37-61		
606-084-00-0	1-(4-methoxy-5-benzofuranyl)-3-phenyl-1,3-propanedione	414-540-3	484-33-3	N; R50-53	N R: 50/53 S: 60-61		
606-085-00-6	(1*R*,4*S*)-2-azabicyclo[2.2.1]hept-5-en-3-one	418-530-1	79200-56-9	Xn; R22 Xi; R41 R43	Xn R: 22-41-43 S: (2-)24-26-37/39		
606-086-00-1	1-(3,3-dimethylcyclohexyl)pent-4-en-1-one	422-330-8	56973-87-6	N; R51-53	N R: 51/53 S: 61		
606-087-00-7	6-ethyl-5-fluoro-4(3*H*)-pyrimidone	422-460-5	137234-87-8	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
606-088-00-2	2,4,4,7-tetramethyl-6-octen-3-one	422-520-0	74338-72-0	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
606-089-00-8	reaction mass of: 1,4-diamino-2-chloro-3-phenoxyanthraquinone; 1,4-diamino-2,3-bis-phenoxyanthraquinone	423-220-2	12223-77-7	R53	R: 53 S: 61		
606-091-00-9	6-chloro-5-(2-chloroethyl)-1,3-dihydroindol-2-one	421-320-0	118289-55-7	N; R50-53	N R: 50/53 S: 60-61		
606-092-00-4	reaction mass of: (*E*)-oxacyclohexadec-12-en-2-one; (*E*)-oxacyclohexadec-13-en-2-one; a) (*Z*)-oxacyclohexadec-(12)-en-2-one and b) (*Z*)-oxacyclohexadec-(13)-en-2-one	422-320-3		N; R50-53	N R: 50/53 S: 60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-001-00-0	formic acid … %	200-579-1	64-18-6	C; R35	C R: 35 S: (1/2-)23-26-45	C; R35: C ≥ 90 % C; R34: 10 % ≤ C < 90 % Xi; R36/38: 2 % ≤ C < 10 %	B
607-002-00-6	acetic acid … %	200-580-7	64-19-7	R10 C; R35	C R: 10-35 S: (1/2-)23-26-45	C; R35: C ≥ 90 % C; R34: 25 % ≤ C < 90 % Xi; R36/38: 10 % ≤ C < 25 %	B
607-003-00-1	chloroacetic acid	201-178-4	79-11-8	T; R25 C; R34 N; R50	T; N R: 25-34-50 S: (1/2-)23-37-45-61		
607-004-00-7	TCA (ISO); trichloroacetic acid	200-927-2	76-03-9	C; R35 N; R50-53	C; N R: 35-50/53 S: (1/2-)26-36/37/39-45-60-61	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
607-005-00-2	TCA-sodium (ISO); sodium trichloroacetate	211-479-2	650-51-1	Xi; R37 N; R50-53	Xi; N R: 37-50/53 S: (2-)46-60-61		
607-006-00-8	oxalic acid	205-634-3	144-62-7	Xn; R21/22	Xn R: 21/22 S: (2-)24/25	Xn; R21/22: C ≥ 5 %	
607-007-00-3	salts of oxalic acid	—	—	Xn; R21/22	Xn R: 21/22 S: (2-)24/25	Xn; R21/22: C ≥ 5 %	A
607-008-00-9	acetic anhydride	203-564-8	108-24-7	R10 Xn; R20/22 C; R34	C R: 10-20/22-34 S: (1/2-)26-36/37/39-45	C; R34: C ≥ 25 % Xi; R37/38-41: 5 % ≤ C < 25 % Xi; R36: 1 % ≤ C < 5 %	
607-009-00-4	phthalic anhydride	201-607-5	85-44-9	Xn; R22 Xi; R37/38-41 R42/43	Xn R: 22-37/38-41-42/43 S: (2-)23-24/25-26-37/39-46		
607-010-00-X	propionic anhydride	204-638-2	123-62-6	C; R34	C R: 34 S: (1/2-)26-45	C; R34: C ≥ 25 % Xi; R36/38: 10 % ≤ C < 25 %	

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-011-00-5	acetyl chloride	200-865-6	75-36-5	F; R11 R14 C; R34	F; C R: 11-14-34 S: (1/2-)9-16-26-45		
607-012-00-0	benzoyl chloride	202-710-8	98-88-4	C; R34	C R: 34 S: (1/2-)26-45		
607-013-00-6	dimethyl carbonate	210-478-4	616-38-6	F; R11	F R: 11 S: (2-)9-16		
607-014-00-1	methyl formate	203-481-7	107-31-3	F+; R12 Xn; R20/22 Xi; R36/37	F+; Xn R: 12-20/22-36/37 S: (2-)9-16-24-26-33		
607-015-00-7	ethyl formate	203-721-0	109-94-4	F; R11 Xn; R20/22 Xi; R36/37	F; Xn R: 11-20/22-36/37 S: (2-)9-16-24-26-33		
607-016-00-2	propyl formate; [1] isopropyl formate [2]	203-798-0 [1] 210-901-2 [2]	110-74-7 [1] 625-55-8 [2]	F; R11 Xi; R36/37 R67	F; Xi R: 11-36/37-67 S: (2-)9-16-24-33		C
607-017-00-8	butyl formate; [1] *tert*-butyl formate; [2] isobutyl formate [3]	209-772-5 [1] 212-105-0 [2] 208-818-1 [3]	592-84-7 [1] 762-75-4 [2] 542-55-2 [3]	F; R11 Xi; R36/37	F; Xi R: 11-36/37 S: (2-)9-16-24-33		C
607-018-00-3	isopentyl formate; [1] pentyl formate; [2] 2-methylbutyl formate [3]	203-769-2 [1] 211-340-6 [2] 252-343-2 [3]	110-45-2 [1] 638-49-3 [2] 35073-27-9 [3]	R10 Xi; R36/37	Xi R: 10-36/37 S: (2-)24		C
607-019-00-9	methyl chloroformate	201-187-3	79-22-1	F; R11 T+; R26 Xn; R21/22 C; R34	F; T+ R: 11-21/22-26-34 S: (1/2-)14-26-28-36/37/39-45-46-63		

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607-020-00-4	ethyl chloroformate	208-778-5	541-41-3	F; R11 T+; R26 Xn; R22 C; R34	F; T+ R: 11-22-26-34 S: (1/2-)9-16-26-28-33-36/37/39-45		
607-021-00-X	methyl acetate	201-185-2	79-20-9	F; R11 Xi; R36 R66 R67	F; Xi R: 11-36-66-67 S: (2-)16-26-29-33		
607-022-00-5	ethyl acetate	205-500-4	141-78-6	F; R11 Xi; R36 R66 R67	F; Xi R: 11-36-66-67 S: (2-)16-26-33		
607-023-00-0	vinyl acetate	203-545-4	108-05-4	F; R11	F R: 11 S: (2-)16-23-29-33		D
607-024-00-6	propyl acetate; [1] isopropyl acetate [2]	203-686-1 [1] 203-561-1 [2]	109-60-4 [1] 108-21-4 [2]	F; R11 Xi; R36 R66 R67	F; Xi R: 11-36-66-67 S: (2-)16-26-29-33		C
607-025-00-1	*n*-butyl acetate	204-658-1	123-86-4	R10 R66 R67	R: 10-66-67 S: (2-)25		
607-026-00-7	*sec*-butyl acetate; [1] isobutyl acetate; [2] *tert*-butyl acetate [3]	203-300-1 [1] 203-745-1 [2] 208-760-7 [3]	105-46-4 [1] 110-19-0 [2] 540-88-5 [3]	F; R11 R66	F R: 11-66 S: (2-)16-23-25-29-33		C
607-027-00-2	methyl propionate	209-060-4	554-12-1	F; R11 Xn; R20	F; Xn R: 11-20 S: (2-)16-24-29-33		
607-028-00-8	ethyl propionate	203-291-4	105-37-3	F; R11	F R: 11 S: (2-)16-23-24-29-33		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-029-00-3	*n*-butyl propionate; [1] *sec*-butyl propionate; [2] *tert*-butyl propionate; [3] iso-butyl propionate [4]	209-669-5 [1] - [2] - [3] 208-746-0 [4]	590-01-2 [1] 591-34-4 [2] 20487-40-5 [3] 540-42-1 [4]	R10	R: 10 S: (2-)		C
607-030-00-9	propyl propionate	203-389-7	106-36-5	R10 Xn; R20	Xn R: 10-20 S: (2-)24		
607-031-00-4	butyl butyrate	203-656-8	109-21-7	R10	R: 10 S: (2-)		C
607-032-00-X	ethyl acrylate	205-438-8	140-88-5	F; R11 Xn; R20/21/22 Xi; R36/37/38 R43	F; Xn R: 11-20/21/22-36/37/38-43 S: (2-)9-16-33-36/37	Xi; 36/37/38: C ≥ 5 %	D
607-033-00-5	*n*-butyl methacrylate	202-615-1	97-88-1	R10 Xi; R36/37/38 R43	Xi R: 10-36/37/38-43 S: (2-)		D
607-034-00-0	methyl acrylate; methyl propenoate	202-500-6	96-33-3	F; R11 Xn; R20/21/22 Xi; R36/37/38 R43	F; Xn R: 11-20/21/22-36/37/38-43 S: (2-)9-25-26-33-36/37-43		D
607-035-00-6	methyl methacrylate; methyl 2-methylprop-2-enoate; methyl 2-methylpropenoate	201-297-1	80-62-6	F; R11 Xi; R37/38 R43	F; Xi R: 11-37/38-43 S: (2-)24-37-46		D
607-036-00-1	2-methoxyethyl acetate; methylglycol acetate	203-772-9	110-49-6	Repr. Cat. 2; R60-61 Xn; R20/21/22	T R: 60-61-20/21/22 S: 53-45		E
607-037-00-7	2-ethoxyethyl acetate; ethylglycol acetate	203-839-2	111-15-9	[$] Repr. Cat. 2; R60-61 Xn; R20/21/22	T R: 60-61-20/21/22 S: 53-45		E

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607-038-00-2	2-butoxyethyl acetate; butylglycol acetate	203-933-3	112-07-2	Xn; R20/21	Xn R: 20/21 S: (2-)24		
607-039-00-8	2,4-D (ISO); 2,4-dichlorophenoxyacetic acid	202-361-1	94-75-7	Xn; R22 Xi; R37-41 R43 R52-53	Xn R: 22-37-41-43-52/53 S: (2-)24/25-26-36/37/39-46-61		
607-040-00-3	salts of 2,4-D	—	—	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)24/25-26-36/37/39-46-61		A
607-041-00-9	2,4,5-T (ISO); 2,4,5-trichlorophenoxy acetic acid	202-273-3	93-76-5	Xn; R22 Xi; R36/37/38 N; R50-53	Xn; N R: 22-36/37/38-50/53 S: (2-)24-60-61		
607-042-00-4	salts and esters of 2,4,5-T; salts and esters of 2,4,5-trichlorophenoxy acetic acid	—	—	Xn; R22 Xi; R36/37/38 N; R50-53	Xn; N R: 22-36/37/38-50/53 S: (2-)24-60-61		A
607-043-00-X	dicamba (ISO); 2,5-dichloro-6-methoxybenzoic acid; 3,6-dichloro-2-methoxybenzoic acid	217-635-6	1918-00-9	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)26-61		
607-044-00-5	3,6-dichloro-*o*-anisic acid, compound with dimethylamine (1:1); [1] potassium 3,6-dichloro-*o*-anisate [2]	218-951-7 [1] 233-002-7 [2]	2300-66-5 [1] 10007-85-9 [2]	Xi; R36 R52-53	Xi R: 36-52/53 S: (2-)26-61		
607-045-00-0	dichlorprop (ISO); 2-(2,4-dichlorophenoxy) propionic acid	204-390-5	120-36-5	Xn; R21/22 Xi; R38-41	Xn R: 21/22-38-41 S: (2-)26-36/37		
607-046-00-6	salts of dichlorprop	—	—	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)13		A
607-047-00-1	fenoprop (ISO); 2-(2,4,5-trichlorophenoxy)propionic acid	202-271-2	93-72-1	Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-50/53 S: (2-)37-60-61		

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607-048-00-7	salts of fenoprop; salts of 2-(2,4,5-trichlorophenoxy)propionic acid	—	—	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61		A
607-049-00-2	mecoprop (ISO); 2-(4-chloro-*o*-tolyloxy) propionic acid; (*RS*)-2-(4-chloro-*o*-tolyloxy)propionic acid; [1] 2-(4-chloro-2-methylphenoxy)propionic acid [2]	230-386-8 [1] 202-264-4 [2]	7085-19-0 [1] 7085-19-0 [2]	Xn; R22 Xi; R38-41 N; R50-53	Xn; N R: 22-38-41-50/53 S: (2-)13-26-37/39-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
607-050-00-8	salts of mecoprop	—	—	Xn; R22 Xi; R38-41 N; R50-53	Xn; N R: 22-38-41-50/53 S: (2-)13-26-37/39-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	A
607-051-00-3	MCPA (ISO); 4-chloro-*o*-tolyloxyacetic acid	202-360-6	94-74-6	Xn; R22 Xi; R38-41	Xn R: 22-38-41 S: (2-)26-37-39		
607-052-00-9	salts and esters of MCPA	—	—	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)13		A
607-053-00-4	MCPB (ISO); 4-(4-chloro-*o*-tolyloxy) butyric acid	202-365-3	94-81-5	N; R50-53	N R: 50/53 S: 60-61		
607-054-00-X	salts and esters of MCPB	—	—	Xn; R22	Xn R: 22 S: (2-)24/25		A
607-055-00-5	endothal-sodium (ISO); disodium 7-oxabicyclo(2,2,1)heptane-2,3-dicarboxylate	204-959-8	129-67-9	T; R25 Xn; R21 Xi; R36/37/38	T R: 21-25-36/37/38 S: (1/2-)36/37/39-45		
607-056-00-0	warfarin (ISO); [1] (*S*)-4-hydroxy-3-(3-oxo-1-phenylbutyl)-2-benzopyrone; [2] (*R*)-4-hydroxy-3-(3-oxo-1-phenylbutyl)-2-benzopyrone [3]	201-377-6 [1] 226-907-3 [2] 226-908-9 [3]	81-81-2 [1] 5543-57-7 [2] 5543-58-8 [3]	Repr. Cat. 1; R61 T; R48/25 R52-53	T R: 61-48/25-52/53 S: 53-45-61		E

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607-057-00-6	coumachlor (ISO); 3-[1-(4-chlorophenyl)-3-oxobutyl]-4-hydroxycoumarin	201-378-1	81-82-3	Xn; R48/22 R52-53	Xn R: 48/22-52/53 S: (2-)37-61		
607-058-00-1	coumafuryl (ISO); fumarin; (*RS*)-3-(1-(2-furyl)-3-oxobutyl)4-hydroxycoumarin; 4-hydroxy-3-[3-oxo-1-(2-furyl) butyl]coumarin	204-195-5	117-52-2	T; R25-48/25 R52-53	T R: 25-48/25-52/53 S: (1/2-)37-45-61		
607-059-00-7	coumatetralyl; 4-hydroxy-3-(1,2,3,4-tetrahydro-1-naphthyl)coumarin	227-424-0	5836-29-3	T+; R27/28 T; R48/24/25 R52-53	T+ R: 27/28-48/24/25-52/53 S: (1/2-)28-36/37-45-61		
607-060-00-2	dicoumarol; 4,4'-dihydroxy-3,3'-methylenebis(2*H*-chromen-2-one)	200-632-9	66-76-2	T; R48/25 Xn; R22 N; R51-53	T; N R: 22-48/25-51/53 S: (1/2-)37-45-61		
607-061-00-8	acrylic acid; prop-2-enoic acid	201-177-9	79-10-7	R10 Xn; R20/21/22 C; R35 N; R50	C; N R: 10-20/21/22-35-50 S: (1/2-)26-36/37/39-45-61	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	D
607-062-00-3	*n*-butyl acrylate	205-480-7	141-32-2	R10 Xi; R36/37/38 R43	Xi R: 10-36/37/38-43 S: (2-)9		D
607-063-00-9	isobutyric acid	201-195-7	79-31-2	Xn; R21/22	Xn R: 21/22 S: (2-)		
607-064-00-4	benzyl chloroformate	207-925-0	501-53-1	C; R34 N; R50-53	C; N R: 34-50/53 S: (1/2-)26-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
607-065-00-X	bromoacetic acid	201-175-8	79-08-3	T; R23/24/25 C; R35 N; R50	T; C; N R: 23/24/25-35-50 S: (1/2-)26-36/37/39-45-61		

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607-066-00-5	dichloroacetic acid	201-207-0	79-43-6	C; R35 N; R50	C; N R: 35-50 S: (1/2-)26-45-61		
607-067-00-0	dichloroacetyl chloride	201-199-9	79-36-7	C; R35 N; R50	C; N R: 35-50 S: (1/2-)9-26-45-61		
607-068-00-6	iodoacetic acid	200-590-1	64-69-7	T; R25 C; R35	T; C R: 25-35 S: (1/2-)22-36/37/39-45		
607-069-00-1	ethyl bromoacetate	203-290-9	105-36-2	T+; R26/27/28	T+ R: 26/27/28 S: (1/2-)7/9-26-45		
607-070-00-7	ethyl chloroacetate	203-294-0	105-39-5	T; R23/24/25 N; R50	T; N R: 23/24/25-50 S: (1/2-)7/9-45-61		
607-071-00-2	ethyl methacrylate	202-597-5	97-63-2	F; R11 Xi; R36/37/38 R43	F; Xi R: 11-36/37/38-43 S: (2-)9-16-29-33		D
607-072-00-8	2-hydroxyethyl acrylate	212-454-9	818-61-1	T; R24 C; R34 R43 N; R50	T; N R: 24-34-43-50 S: (1/2-)26-36/39-45-61	T; R24: C ≥ 2 % Xn; R21: 0,2 % ≤ C < 2 % R43: C ≥ 0,2 %	D
607-073-00-3	4-CPA (ISO); 4-chlorophenoxyacetic acid	204-581-3	122-88-3	Xn; R22	Xn R: 22 S: (2-)		
607-074-00-9	chlorfenac (ISO); 2,3,6-trichlorophenylacetic acid	201-599-3	85-34-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)36-61		
607-075-00-4	chlorfenprop-methyl; methyl 2-chloro-3-(4-chlorophenyl)propionate	238-413-5	14437-17-3	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		

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607-076-00-X	dodine (ISO); dodecylguanidinium acetate	219-459-5	2439-10-3	Xn; R22 Xi; R36/38 N; R50-53	Xn; N R: 22-36/38-50/53 S: (2-)26-60-61		
607-077-00-5	erbon (ISO); 2-(2,4,5-trichlorophenoxy)ethyl 2,2-dichloropropionate	—	136-25-4	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
607-078-00-0	fluenetil (ISO); 2-fluoroethyl biphenyl-4-ylacetate	—	4301-50-2	T+; R27/28	T+ R: 27/28 S: (1/2-)28-36/37-45		
607-079-00-6	kelevan (ISO); ethyl 5-(perchloro-5-hydroxypentacyclo[5,3,0,$0^{2,6}$,$0^{3,9}$,$0^{4,8}$]decan-5-yl)-4-oxopentanoate; ethyl 5-(1,2,3,5,6,7,8,9,10,10-decachloro-4-hydroxypentacyclo(5,2,1,$0^{2,6}$,$0^{3,9}$,$0^{5,8}$)dec-4-yl)-4-oxovalerate	—	4234-79-1	T; R24 Xn; R22 N; R51-53	T; N R: 22-24-51/53 S: (1/2-)36/37-45-61		
607-080-00-1	chloroacetyl chloride	201-171-6	79-04-9	R14 R29 T; R23/24/25-48/23 C; R35 N; R50	T; C; N R: 14-23/24/25-29-35-48/23-50 S: (1/2-)7/8-9-26-36/37/39-45-61		
607-081-00-7	fluoroacetic acid	205-631-7	144-49-0	T+; R28 N; R50	T+; N R: 28-50 S: (1/2-)20-22-26-45-61		
607-082-00-2	fluoroacetates, soluble	—	—	T+; R28 N; R50	T+; N R: 28-50 S: (1/2-)20-22-26-45-61		A
607-083-00-8	2,4-DB (ISO); 4-(2,4-dichlorophenoxy)butyric acid	202-366-9	94-82-6	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)25-29-46-61		
607-084-00-3	salts of 2,4-DB	—	—	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)26-29-39-46-61		A

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607-085-00-9	benzyl benzoate	204-402-9	120-51-4	Xn; R22	Xn R: 22 S: (2-)25		
607-086-00-4	diallyl phthalate	205-016-3	131-17-9	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24/25-60-61		
607-088-00-5	methacrylic acid; 2-methylpropenoic acid	201-204-4	79-41-4	Xn; R21/22 C; R35	C R: 21/22-35 S: (1/2-)26-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	D
607-089-00-0	propionic acid … %	201-176-3	79-09-4	C; R34	C R: 34 S: (1/2-)23-36-45	C; R34: C ≥ 25 % Xi; R36/37/38: 10 % ≤ C < 25 %	B
607-090-00-6	thioglycolic acid	200-677-4	68-11-1	T; R23/24/25 C; R34	T R: 23/24/25-34 S: (1/2-)25-27-28-45	T; R23/24/25: C ≥ 2 % Xn; R20/21/22: 0,2 % ≤ C < 2 %	
607-091-00-1	trifluoroacetic acid . . . %	200-929-3	76-05-1	Xn; R20 C; R35 R52-53	C R: 20-35-52/53 S: (1/2-)9-26-27-28-45-61	Xn; R20: C ≥ 10 %	B
607-092-00-7	methyl lactate; [1] methyl (±)-lactate; [2] methyl (*R*)-lactate; [3] methyl (*S*)-(-)-lactate [4]	208-930-0 [1] 218-449-8 [2] 241-420-6 [3] 248-704-9 [4]	547-64-8 [1] 2155-30-8 [2] 17392-83-5 [3] 27871-49-4 [4]	R10 Xi; R36/37	Xi R: 10-36/37 S: (2-)24		C
607-093-00-2	propionyl chloride	201-170-0	79-03-8	F; R11 R14 C; R34	F; C R: 11-14-34 S: (1/2-)9-16-26-45		B D
607-094-00-8	peracetic acid . . . %	201-186-8	79-21-0	R10 O; R7 Xn; R20/21/22 C; R35 N; R50	O; C; N R: 7-10-20/21/22-35-50 S: (1/2-)3/7-14-36/37/39-45-61	Xn; R20/21/22: C ≥ 10 % C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	B D

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607-095-00-3	maleic acid	203-742-5	110-16-7	Xn; R22 Xi; R36/37/38	Xn R: 22-36/37/38 S: (2-)26-28-37		
607-096-00-9	maleic anhydride	203-571-6	108-31-6	Xn; R22 C; R34 R42/43	C R: 22-34-42/43 S: (2-)22-26-36/37/39-45		
607-097-00-4	benzene-1,2,4-tricarboxylic acid 1,2-anhydride; trimellitic anhydride	209-008-0	552-30-7	Xi; R37-41 R42/43	Xn R: 37-41-42/43 S: (2-)22-26-36/37/39		
607-098-00-X	benzene-1,2:4,5-tetracarboxylic dianhydride; benzene-1,2:4,5-tetracarboxylic dianhydride; pyromellitic dianhydride	201-898-9	89-32-7	Xi; R41 R42/43	Xn R: 41-42/43 S: (2-)22-24-26-37/39		
607-099-00-5	1,2,3,6-tetrahydrophthalic anhydride; [1] *cis*-1,2,3,6-tetrahydrophthalic anhydride; [2] 3,4,5,6-tetrahydrophthalic anhydride; [3] tetrahydrophthalic anhydride [4]	201-605-4 [1] 213-308-7 [2] 219-374-3 [3] 247-570-9 [4]	85-43-8 [1] 935-79-5 [2] 2426-02-0 [3] 26266-63-7 [4]	Xi; R41 R42/43 R52-53	Xn R: 41-42/43-52/53 S: (2-)22-24-26-37/39-61		C
607-100-00-9	benzophenone-3,3',4,4'-tetracarboxylic dianhydride; 4,4'-carbonyldi(phthalic anhydride)	219-348-1	2421-28-5	Xi; R36/37	Xi R: 36/37 S: (2-)25	Xi; R36/37: C ≥ 1 %	
607-101-00-4	1,4,5,6,7,7-hexachlorobicyclo [2,2,1]hept-5-ene-2,3-dicarboxylic anhydride chlorendic anhydride	204-077-3	115-27-5	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)25	Xi; R36/37/38: C ≥ 1 %	
607-102-00-X	cyclohexane-1,2-dicarboxylic anhydride; [1] *cis*-cyclohexane-1,2-dicarboxylic anhydride; [2] *trans*-cyclohexane-1,2-dicarboxylic anhydride [3]	201-604-9 [1] 236-086-3 [2] 238-009-9 [3]	85-42-7 [1] 13149-00-3 [2] 14166-21-3 [3]	Xi; R41 R42/43	Xn R: 41-42/43 S: (2-)23-24-26-37/39		C
607-103-00-5	succinic anhydride	203-570-0	108-30-5	Xi; R36/37	Xi R: 36/37 S: (2-)25	Xi; R36/37: C ≥ 1 %	

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607-104-00-0	cyclopentane-1,2,3,4-tetracarboxylic dianhydride	227-964-7	6053-68-5	Xi; R36/37	Xi R: 36/37 S: (2-)25	Xi; R36/37: C ≥ 1 %	
607-105-00-6	8,9,10-trinorborn-5-ene-2,3-dicarboxylic anhydride; [1] 1,2,3,6-tetrahydro-3,6-methanophthalic anhydride; [2] (1α,2α,3β,6β)-1,2,3,6-tetrahydro-3,6-methanophthalic anhydride [3]	204-957-7 [1] 212-557-9 [2] 220-384-5 [3]	129-64-6 [1] 826-62-0 [2] 2746-19-2 [3]	Xi; R41 R42/43	Xn R: 41-42/43 S: (2-)22-24-26-37/39		C
607-106-00-1	8,9-dinorborn-5-ene-2,3-dicarboxylic anhydride	—	123748-85-6	Xn; R22 Xi; R36/37/38 R42	Xn R: 22-36/37/38-42 S: (2-)39	Xi; R36/37/38: C ≥ 10 %	C
607-107-00-7	2-ethylhexyl acrylate	203-080-7	103-11-7	Xi; R37/38 R43	Xi R: 37/38-43 S: (2-)36/37-46		D
607-108-00-2	2-hydroxy-1-methylethylacrylate; [1] 2-hydroxypropylacrylate; [2] acrylic acid, monoester with propane-1,2-diol [3]	220-852-9 [1] 213-663-8 [2] 247-118-0 [3]	2918-23-2 [1] 999-61-1 [2] 25584-83-2 [3]	T; R23/24/25 C; R34 R43	T R: 23/24/25-34-43 S: (1/2-)26-36/37/39-45	T; R23/24/25: C ≥ 2 % Xn; R20/21/22: 0,2 % ≤ C < 2 % R43: C ≥ 0,2 %	C D
607-109-00-8	hexamethylene diacrylate; hexane-1,6-diol diacrylate	235-921-9	13048-33-4	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)39		D
607-110-00-3	pentaerythritol triacrylate	222-540-8	3524-68-3	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)39		D
607-111-00-9	2,2-bis(acryloyloxymethyl)butyl acrylate; trimethylolpropane triacrylate	239-701-3	15625-89-5	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)39		D
607-112-00-4	2,2-dimethyltrimethylene diacrylate; neopentyl glycol diacrylate	218-741-5	2223-82-7	T; R24 Xi; R36/38 R43	T R: 24-36/38-43 S: (1/2-)28-39-45	T; R24: C ≥ 5 % Xn; R21: 0,2 % ≤ C < 5 %	D

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607-113-00-X	isobutyl methacrylate	202-613-0	97-86-9	R10 Xi; R36/37/38 R43 N; R50	Xi; N R: 10-36/37/38-43-50 S: (2-)24-37-61		D
607-114-00-5	ethylene dimethacrylate	202-617-2	97-90-5	Xi; R37 R43	Xi R: 37-43 S: (2-)24-37	Xi; R37: C ≥ 10 %	D
607-115-00-0	isobutyl acrylate	203-417-8	106-63-8	R10 Xn; R20/21 Xi; R38 R43	Xn R: 10-20/21-38-43 S: (2-)9-24-37	Xi; R38: C ≥ 10 %	D
607-116-00-6	cyclohexyl acrylate	221-319-3	3066-71-5	Xi; R37/38 N; R51-53	Xi; N R: 37/38-51/53 S: (2-)61	Xi; R37/38: C ≥ 10 %	D
607-117-00-1	2,3-epoxypropyl acrylate; glycidyl acrylate	203-440-3	106-90-1	T; R23/24/25 C; R34 R43	T R: 23/24/25-34-43 S: (1/2-)26-36/37/39-45	T; R23/24/25: C ≥ 2 % Xn; R20/21/22: 0,2 % ≤ C < 2 % R43: C ≥ 0,2 %	D
607-118-00-7	1-methyltrimethylene diacrylate; 1,3-butylene glycol diacrylate	243-105-9	19485-03-1	Xn; R21 C; R34 R43	C R: 21-34-43 S: (1/2-)26-36/37/39-45		D
607-119-00-2	tetramethylene diacrylate; 1,4-butyleneglycol diacrylate	213-979-6	1070-70-8	Xn; R21 C; R34 R43	C R: 21-34-43 S: (1/2-)26-36/37/39-45		D
607-120-00-8	2,2'-oxydiethyl diacrylate; diethylene glycol diacrylate	223-791-6	4074-88-8	T; R24 Xi; R36/38 R43	T R: 24-36/38-43 S: (1/2-)28-39-45	T; R24: C ≥ 2 % Xn; R21: 0,2 % ≤ C < 2 % R43: C ≥ 0,2 %	D
607-121-00-3	8,9,10-trinorborn-2-yl acrylate	—	10027-06-2	Xn; R21 Xi; R38 R43	Xn R: 21-38-43 S: (2-)28	Xi; R38: C ≥ 10 %	D

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607-122-00-9	pentaerythritol tetraacrylate	225-644-1	4986-89-4	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)26-39		D
607-123-00-4	2,3-epoxypropyl methacrylate; glycidyl methacrylate	203-441-9	106-91-2	Xn; R20/21/22 Xi; R36/38 R43	Xn R: 20/21/22-36/38-43 S: (2-)26-28	Xi; R36/38: C ≥ 10 %	D
607-124-00-X	2-hydroxyethyl methacrylate	212-782-2	868-77-9	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)26-28		D
607-125-00-5	2-hydroxypropyl methacrylate; [1] 3-hydroxypropyl methacrylate [2]	213-090-3 [1] 220-426-2 [2]	923-26-2 [1] 2761-09-3 [2]	Xi; R36 R43	Xi R: 36-43 S: (2-)24/25-26-37/39		C D
607-126-00-0	2,2'-(ethylenedioxy)diethyl diacrylate; triethylene glycol diacrylate	216-853-9	1680-21-3	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)26-28		D
607-127-00-6	2-diethylaminoethyl methacrylate	203-275-7	105-16-8	Xn; R20 Xi; R36/38 R43	Xn R: 20-36/38-43 S: (2-)26	Xi; R36/38: C ≥ 10 %	D
607-128-00-1	2-*tert*-butylaminoethyl methacrylate	223-228-4	3775-90-4	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)26		D
607-129-00-7	ethyl lactate; ethyl DL-lactate; [1] ethyl (*S*)-2-hydroxypropionate; ethyl L-lactate; ethyl-(*S*)-lactate [2]	202-598-0 [1] 211-694-1 [2]	97-64-3 [1] 687-47-8 [2]	R10 Xi; R37-41	Xi R: 10-37-41 S: (2-)24-26-39		C

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-130-00-2	pentyl acetate; [1] isopentyl acetate; [2] 1-methylbutyl acetate; [3] 2-methylbutyl acetat; [4] 2(or 3)-methylbutyl acetate [5]	211-047-3 [1] 204-662-3 [2] 210-946-8 [3] 210-843-8 [4] 282-263-3 [5]	628-63-7 [1] 123-92-2 [2] 626-38-0 [3] 624-41-9 [4] 84145-37-9 [5]	R10 R66	R: 10-66 S: (2-)23-25		C
607-131-00-8	isopentyl propionate; [1] pentyl propionate; [2] 2-methylbutyl propionate [3]	203-322-1 [1] 210-852-7 [2] 219-449-0 [3]	105-68-0 [1] 624-54-4 [2] 2438-20-2 [3]	R10	R: 10 S: (2-)23-24		C
607-132-00-3	2-dimethylaminoethyl methacrylate	220-688-8	2867-47-2	Xn; R21/22 Xi; R36/38 R43	Xn R: 21/22-36/38-43 S: (2-)26-28	Xi; R36/38: C ≥ 10 %	D
607-133-00-9	monoalkyl or monoaryl or monoalkylaryl esters of acrylic acid with the exception of those specified elsewhere in this Annex	—	—	Xi; R36/37/38 N; R51-53	Xi; N R: 36/37/38-51/53 S: (2-)26-28-61	Xi; R36/37/38: C ≥ 10 %	A
607-134-00-4	monoalkyl or monoaryl or monoalkyaryl esters of methacrylic acid with the exception of those specified elsewhere in this Annex	—	—	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)26-28	Xi; R36/37/38: C ≥ 10 %	A
607-135-00-X	butyric acid	203-532-3	107-92-6	C; R34	C R: 34 S: (1/2-)26-36-45		
607-136-00-5	butyryl chloride	205-498-5	141-75-3	F; R11 C; R34	F; C R: 11-34 S: (1/2-)16-23-26-36-45		
607-137-00-0	methyl acetoacetate	203-299-8	105-45-3	Xi; R36	Xi R: 36 S: (2-)26		

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607-138-00-6	butyl chloroformate; chloroformic acid butyl ester	209-750-5	592-34-7	R10 T; R23 C; R34	T R: 10-23-34 S: (1/2-)26-36-45		
607-139-00-1	2-chloropropionic acid	209-952-3	598-78-7	Xn; R22 C; R35	C R: 22-35 S: (1/2-)23-26-28-36-45		
607-140-00-7	isobutyryl chloride	201-194-1	79-30-1	F; R11 C; R35	F; C R: 11-35 S: (1/2-)16-23-26-36-45		
607-141-00-2	oxydiethylene bis(chloroformate)	203-430-9	106-75-2	Xn; R22 Xi; R38-41 N; R51-53	Xn; N R: 22-38-41-51/53 S: (2-)23-26-61		
607-142-00-8	propyl chloroformate; chloroformic acid propylester; *n*-propyl chloroformate	203-687-7	109-61-5	R10 [$] T; R23 C; R34	T R: 10-23-34 S: (1/2-)26-36-45		
607-143-00-3	valeric acid	203-677-2	109-52-4	C; R34 R52-53	C R: 34-52/53 S: (1/2-)26-36-45-61		
607-144-00-9	adipic acid	204-673-3	124-04-9	Xi; R36	Xi R: 36 S: (2-)		
607-145-00-4	methanesulphonic acid	200-898-6	75-75-2	C; R34	C R: 34 S: (1/2-)26-36-45		
607-146-00-X	fumaric acid	203-743-0	110-17-8	Xi; R36	Xi R: 36 S: (2-)26		
607-147-00-5	oxalic acid diethylester; diethyl oxalate	202-464-1	95-92-1	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)23		

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607-148-00-0	guanidinium chloride; guanadine hydrochloride	200-002-3	50-01-1	Xn; R22 Xi; R36/38	Xn R: 22-36/38 S: (2-)22		
607-149-00-6	urethane (INN); ethyl carbamate	200-123-1	51-79-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		
607-150-00-1	endothal (ISO); 7-oxabicyclo(2,2,1)heptane-2,3-dicarboxylic acid	205-660-5	145-73-3	T; R25 Xn; R21 Xi; R36/37/38	T R: 21-25-36/37/38 S: (1/2-)36/37/39-45		
607-151-00-7	propargite (ISO); 2-(4-*tert*-butylphenoxy) cyclohexyl prop-2-ynyl sulphite	219-006-1	2312-35-8	Carc. Cat. 3; R40 T; R23 Xi; R38-41 N; R50-53	T; N R: 23-38-40-41-50/53 S: (1/2-)26-36/37/39-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
607-152-00-2	2,3,6-TBA (ISO); 2,3,6-trichlorobenzoic acid	200-026-4	50-31-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
607-153-00-8	benazolin (ISO); 4-chloro-2,3-dihydro-2-oxo-1,3-benzothiazol-3-ylacetic acid	223-297-0	3813-05-6	Xi; R36/38 R52-53	Xi R: 36/38-52/53 S: (2-)22-61		
607-154-00-3	ethyl *N*-benzoyl-*N*-(3,4-dichlorophenyl)-DL-alaninate; benzoylprop-ethyl (ISO)	244-845-5	22212-55-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24-60-61		
607-155-00-9	3-(3-amino-5-(1-methylguanidino)-1-oxopentylamino-6-(4-amino-2-oxo-2,3-dihydro-pyrimidin-1-yl)-2,3-dihydro-(6*H*)-pyran-2-carboxylic acid; blasticidin-s	—	2079-00-7	T+; R28	T+ R: 28 S: (1/2-)24/25-36/37-45		
607-156-00-4	chlorfenson (ISO); 4-chlorophenyl 4-chlorobenzenesulfonate	201-270-4	80-33-1	Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-50/53 S: (2-)37-60-61		
607-157-00-X	3-(3-biphenyl-4-yl-1,2,3,4-tetrahydro-1-naphthyl)-4-hydroxy-coumarin; difenacoum	259-978-4	56073-07-5	T+; R28 T; R48/25 N; R50-53	T+; N R: 28-48/25-50/53 S: (1/2-)36/37-45-60-61		

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607-158-00-5	sodium salt of chloroacetic acid; sodium chloroacetate	223-498-3	3926-62-3	T; R25 Xi; R38 N; R50	T; N R: 25-38-50 S: (1/2-)22-37-45-61		
607-159-00-0	chlorobenzilate (ISO); ethyl 2,2-di(4-chlorophenyl)-2-hydroxyacetate; ethyl 4,4'-dichlorobenzilate	208-110-2	510-15-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
607-160-00-6	isobutyl 2-(4-(4-chlorophenoxy)phenoxy)propionate; clofop-isobutyl (ISO)	—	51337-71-4	Xn; R22	Xn R: 22 S: (2-)		
607-161-00-1	diethanolamine salt of 4-CPA	—	—	Xn; R22	Xn R: 22 S: (2-)		
607-162-00-7	2,2-dichloropropionic acid; dalapon	200-923-0	75-99-0	Xn; R22 Xi; R38-41 R52-53	Xn R: 22-38-41-52/53 S: (2-)26-39-61		
607-163-00-2	3-acetyl-6-methyl-2*H*-pyran-2,4(3*H*)-dione; dehydracetic acid	208-293-9	520-45-6	Xn; R22	Xn R: 22 S: (2-)		
607-164-00-8	sodium 1-(3,4-dihydro-6-methyl-2,4-dioxo-2*H*-pyran-3-ylidene)ethonolate; sodium dehydracetate	224-580-1	4418-26-2	Xn; R22	Xn R: 22 S: (2-)		
607-165-00-3	diclofop-methyl (ISO); methyl 2-(4-(2,4-dichlorophenoxy)phenoxy)propionate; methyl (*RS*)-2-[4-(2,4-dichlorophenoxy)phenoxy]propionate;	257-141-8	51338-27-3	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
607-166-00-9	medinoterb acetate (ISO); 6-*tert*-butyl-3-methyl-2,4-dinitrophenyl acetate	219-634-6	2487-01-6	T; R25 Xn; R21	T R: 21-25 S: (1/2-)36/37-45		
607-167-00-4	sodium 3-chloroacrylate	—	4312-97-4	Xn; R21/22	Xn R: 21/22 S: (2-)36/37		

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607-168-00-X	dipropyl 6,7-methylenedioxy-1,2,3,4-tetrahydro-3-methyl-naphthalene-1,2-dicarboxylate; propylisome	—	83-59-0	T; R24 Xn; R22 N; R50-53	T; N R: 22-24-50/53 S: (1/2-)36/37-45-60-61		
607-169-00-5	sodium fluoroacetate	200-548-2	62-74-8	T+; R26/27/28 N; R50	T+; N R: 26/27/28-50 S: (1/2-)13-22-36/37-45-61		
607-170-00-0	bis(1,2,3-trithiacyclohexyldimethylammonium) oxalate; thiocyclam-oxalate	250-859-2	31895-22-4	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-46-60-61		
607-172-00-1	4-hydroxy-3-(3-(4'-bromo-4-biphenylyl)-1,2,3,4-tetrahydro-1-naphthyl)coumarin; brodifacoum	259-980-5	56073-10-0	T+; R27/28 T; R48/24/25 N; R50-53	T+; N R: 27/28-48/24/25-50/53 S: (1/2-)36/37-45-60-61		
607-173-00-7	dimethyl (3-methyl-4-(5-nitro-3-ethoxycarbonyl-2-thienyl) azo)phenylnitrilodipropionate	400-460-6	—	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
607-174-00-2	reaction mass of dodecyl 3-(2,2,4,4-tetramethyl-21-oxo-7-oxa-3,20-diazadispiro(5,1,11,2)henicosan-20-yl)propionate and tetradecyl 3-(2,2,4,4-tetramethyl-21-oxo-7-oxa-3,20-diazadispiro(5,1,11,2)henicosan-20-yl)propionate	400-580-9	—	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)28-61		
607-175-00-8	methyl 2-(2-nitrobenzylidene)acetoacetate	400-650-9	39562-27-1	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-176-00-3	reaction mass of α-3-(3-(2*H*-benzotriazol-2-yl)-5-*tert*-butyl-4-hydroxyphenyl)propionyl-ω-hydroxypoly(oxyethylene) and α-3-(3-(2*H*-benzotriazol-2-yl)-5-*tert*-butyl-4-hydroxyphenyl) propionyl-ω-3-(3-(2*H*-benzotriazol-2-yl)-5-*tert*-butyl-4-hydroxyphenyl)propionyloxypoly(oxyethylene)	400-830-7	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)36/37-61		
607-177-00-9	methyl 2-(3-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)3-methylureidosulphonyl)benzoate	401-190-1	101200-48-0	R43	Xi R: 43 S: (2-)22-24-37		
607-178-00-4	methyl α-((4,6-dimethoxypyrimidin-2-yl)ureidosulphonyl)-*o*-toluate	401-340-6	83055-99-6	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		

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607-179-00-X	(benzothiazol-2-ylthio)succinic acid	401-450-4	95154-01-1	R43	Xi R: 43 S: (2-)24-37		
607-180-00-5	potassium 2-hydroxycarbazole-1-carboxylate	401-630-2	96566-70-0	Xn; R22 Xi; R36/37 R52-53	Xn R: 22-36/37-52/53 S: (2-)22-26-61		
607-181-00-0	3,5-dichloro-2,4-difluorobenzoyl fluoride	401-800-6	101513-70-6	T; R23 C; R34 Xn; R22 R29 R43 R52-53	T; C R: 22-23-29-34-43-52/53 S: (1/2-)26-36/37/39-45-61		
607-182-00-6	methyl 3-sulphamoyl-2-thenoate	402-050-2	—	R43	Xi R: 43 S: (2-)24-37		
607-183-00-1	zinc 2-hydroxy-5-C_{13-18}alkylbenzoate	402-280-3	—	Xi; R36/38 N; R51-53	Xi; N R: 36/38-51/53 S: (2-)26-61		
607-184-00-7	*S*-(3-trimethoxysilyl)propyl 19-isocyanato-11-(6-isocyanatohexyl)-10,12-dioxo-2,9,11,13-tetraazanonadecanethioate	402-290-8	85702-90-5	R10 R42/43	Xn R: 10-42/43 S: (2-)23-24-37		
607-185-00-2	ethyl *trans*-3-dimethylaminoacrylate	402-650-4	1117-37-9	R43	Xi R: 43 S: (2-)24-37		
607-186-00-8	quinclorac (ISO); 3,7-dichloroquinoline-8-carboxylic acid	402-780-1	84087-01-4	R43	Xi R: 43 S: (2-)24-37		
607-187-00-3	bis(2,2,6,6-tetramethyl-4-piperidyl) succinate	402-940-0	62782-03-0	Xi; R36 R52-53	Xi R: 36-52/53 S: (2-)26-61		
607-188-00-9	hydrogen sodium *N*-carboxylatoethyl-*N*-octadec-9-enylmaleamate	402-970-4	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-189-00-4	trimethylenediaminetetraacetic acid	400-400-9	1939-36-2	Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-50/53 S: (2-)22-26-39-60-61		
607-190-00-X	methyl acrylamidomethoxyacetate (containing ≥ 0,1 % acrylamid)	401-890-7	77402-03-0	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Xn; R22 Xi; R36	T R: 45-46-22-36 S: 53-45		E
607-191-00-5	isobutyl 3,4-epoxybutyrate	401-920-9	100181-71-3	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-28-36/37-60-61		
607-192-00-0	disodium *N*-carboxymethyl-*N*-(2-(2-hydroxyethoxy)ethyl)glycinate	402-360-8	92511-22-3	Xi; R41	Xi R: 41 S: (2-)26-39		
607-194-00-1	propylene carbonate	203-572-1	108-32-7	Xi; R36	Xi R: 36 S: (2-)		
607-195-00-7	2-methoxy-1-methylethyl acetate	203-603-9	108-65-6	R10 Xi; R36	Xi R: 10-36 S: (2-)25		
607-196-00-2	heptanoic acid	203-838-7	111-14-8	C; R34	C R: 34 S: (1/2-)26-28-36/37/39-45		
607-197-00-8	nonanoic acid	203-931-2	112-05-0	C; R34	C R: 34 S: (1/2-)26-28-36/37/39-45		
607-198-00-3	propyl 3,4,5-trihydroxybenzoate	204-498-2	121-79-9	Xn; R22 R43	Xn R: 22-43 S: (2-)24-37		

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607-199-00-9	octyl 3,4,5-trihydroxybenzoate	213-853-0	1034-01-1	Xn; R22 R43	Xn R: 22-43 S: (2-)24-37		
607-200-00-2	dodecyl 3,4,5-trihydroxybenzoate	214-620-6	1166-52-5	R43	Xi R: 43 S: (2-)24-37		
607-201-00-8	thiocarbonyl chloride	207-341-6	463-71-8	T; R23 Xn; R22 Xi; R36/37/38	T R: 22-23-36/37/38 S: (1/2-)7-9-36/37-45		
607-203-00-9	2-ethylhexyl[[[3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl]methyl]thio]acetate	279-452-8	80387-97-9	Repr. Cat. 2; R61 R43 R52-53	T R: 61-43-52/53 S: 53-45-61		
607-204-00-4	(chlorophenyl)(chlorotolyl)methane, mixed isomers	400-140-6	—	N; R50-53	N R: 50/53 S: 60-61		
607-205-00-X	methyl chloroacetate	202-501-1	96-34-4	R10 T; R23/25 Xi; R37/38-41	T R: 10-23/25-37/38-41 S: (1/2-)26-37/39-45		
607-206-00-5	isopropyl chloroacetate	203-301-7	105-48-6	R10 T; R25 Xi; R36/37/38	T R: 10-25-36/37/38 S: (1/2-)26-37/39-45		
607-207-00-0	haloxyfop-etotyl (ISO) 2-ethoxyethyl 2-(4-(3-chloro-5-trifluoromethyl-2-pyridyloxy)phenoxy)propionate; haloxyfop-(2-ethoxyethyl)	402-560-5	87237-48-7	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-36-60-61		
607-208-00-6	4,8,12-trimethyltrideca-3,7,11-trienoic acid, mixed isomers	403-000-2	91853-67-7	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)37/39-60-61		
607-209-00-1	reaction mass of O,O'-diisopropyl (pentathio)dithioformate and O,O'-diisopropyl (trithio)dithioformate and O,O'-diisopropyl (tetrathio)dithioformate	403-030-6	—	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)36/37-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-210-00-7	methyl acrylamidoglycolate (containing ≥ 0,1 % acrylamide)	403-230-3	77402-05-2	Carc. Cat. 2; R45 Muta. Cat. 2; R46 C; R34 R43	T R: 45-46-34-43 S: 53-45		
607-211-00-2	methyl 3-(3-*tert*-butyl-4-hydroxy-5-methylphenyl)propionate	403-270-1	6386-39-6	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)36-61		
607-212-00-8	poly(oxypropylenecarbonyl-co-oxy(ethylethylene)carbonyl), containing 27 % hydroxyvalerate	403-300-3	—	R43	Xi R: 43 S: (2-)24-37		
607-213-00-3	ethyl 3,3-bis(*tert*-pentylperoxy)butyrate	403-320-2	67567-23-1	E; R2 [$] O; R7 R10 N; R51-53	E; N R: 2-7-10-51/53 S: (2-)3/7-14-33-36/37/39-61		
607-214-00-9	*N*,*N*-hydrazinodiacetic acid	403-510-5	19247-05-3	T; R25 Xn; R48/22 R43 R52-53	T R: 25-43-48/22-52/53 S: (1/2-)26-36/37/39-45-61		
607-215-00-4	3-(3-*tert*-butyl-4-hydroxyphenyl)propionic acid	403-920-4	107551-67-7	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)25-26-36		
607-216-00-X	glutamic acid, reaction products with *N*-(C_{12-14}alkyl)propylenediamine	403-950-8	—	T+; R26 Xn; R22 C; R34 N; R50-53	T+; N R: 22-26-34-50/53 S: (1/2-)26-36/37/39-38-45-60-61		
607-217-00-5	2-ethoxyethyl 2-(4-(2,6-dihydro-2,6-dioxo-7-phenyl-1,5-dioxaindacen-3-yl)phenoxy)acetate	403-960-2	—	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
607-218-00-0	dichlorprop-P (ISO); (+)-*R*-2-(2,4-dichlorophenoxy)propionic acid	403-980-1	15165-67-0	Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43 S: (2-)24-26-37/39		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-219-00-6	bis(2-ethylhexyl) dithiodiacetate	404-510-8	62268-47-7	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)24/25-37-61		
607-221-00-7	6-docosyloxy-1-hydroxy-4-(1-(4-hydroxy-3-methylphenanthren-1-yl)-3-oxo-2-oxaphenalen-1-yl)naphthalene-2-carboxylic acid	404-550-6	—	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
607-222-00-2	6-(2,3-dimethylmaleimido)hexyl methacrylate	404-870-6	63740-41-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-223-00-8	transfluthrin (ISO); 2,3,5,6-tetrafluorobenzyl *trans*-2-(2,2-dichlorovinyl)-3,3-dimethylcyclopropanecarboxylate	405-060-5	118712-89-3	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)36/37-60-61		
607-224-00-3	methyl 2-(3-nitrobenzylidene)acetoacetate	405-270-7	39562-17-9	Xi; R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
607-225-00-9	3-azidosulfonylbenzoic acid	405-310-3	15980-11-7	E; R2 Xn; R48/22 Xi; R41 R43	E; Xn R: 2-41-43-48/22 S: (2-)22-26-35-36/37/39		
607-226-00-4	reaction mass of 2-acryloyloxyethyl hydrogen cyclohexane-1,2-dicarboxylate and 2-methacryloyloxyethyl hydrogen cyclohexane-1,2-dicarboxylate	405-360-6	—	Xi; R38-41 R43 R52-53	Xi R: 38-41-43-52/53 S: (2-)24-26-37/39-61		
607-227-00-X	potassium 2-amino-2-methylpropionate octahydrate	405-560-3	120447-91-8	Xn; R22 C; R35	C R: 22-35 S: (1/2-)26-28-36/37/39-45		
607-228-00-5	bis(2-methoxyethyl) phthalate	204-212-6	117-82-8	Repr. Cat. 2; R61 Repr. Cat. 3; R62	T R: 61-62 S: 53-45		
607-229-00-0	diethylcarbamoyl chloride	201-798-5	88-10-8	Carc. Cat. 3; R40 Xn; R20/22 Xi; R36/37/38	Xn R: 20/22-36/37/38-40 S: (2-)26-36/37		

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607-230-00-6	2-ethylhexanoic acid	205-743-6	149-57-5	Repr. Cat. 3; R63	Xn R: 63 S: (2-)36/37		
607-231-00-1	3,6-dichloropyridine-2-carboxylic acid; clopyralid	216-935-4	1702-17-6	Xi; R41 N; R51-53	X; N R: 41-51/53 S: (2-)26-39-61		
607-232-00-7	pyridate (ISO); *O*-(6-chloro-3-phenylpyridazin-4-yl) *S*-octyl thiocarbonate	259-686-7	55512-33-9	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
607-233-00-2	hexyl acrylate	219-698-5	2499-95-8	Xi; R36/37/38 R43 N; R51-53	Xi; N R: 36/37/38-43-51/53 S: (2-)24-26-37-61		
607-234-00-8	flurenol (ISO); 9-hydroxy-9*H*-fluorene-9-carboxylic acid	207-397-1	467-69-6	N; R51-53	N R: 51/53 S: 61		
607-235-00-3	mecrilate; methyl 2-cyanoacrylate	205-275-2	137-05-3	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)23-24/25-26	Xi; R36/37/38: C ≥ 10 %	
607-236-00-9	ethyl 2-cyanoacrylate	230-391-5	7085-85-0	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)23-24/25-26	Xi; R36/37/38: C ≥ 10 %	
607-237-00-4	benzyl 2-chloro-4-(trifluoromethyl)thiazole-5-carboxylate; flurazole	276-942-3	72850-64-7	N; R51-53	N R: 51/53 S: 61		
607-238-00-X	tau-fluvalinate (ISO); cyano-(3-phenoxyphenyl)methyl *N*-[2-chloro-4-(trifluoromethyl)phenyl]-D-valinate	—	102851-06-9	Xn; R22 Xi; R38 N; R50-53	Xn; N R: 22-38-50/53 S: (2-)24-59-61		
607-239-00-5	fenpropathrin (ISO); α-cyano-3-phenoxybenzyl 2,2,3,3-tetramethylcyclopropanecarboxylate;	254-485-0	39515-41-8	T+; R26 T; R25 Xn; R21 N; R50-53	T+; N R: 21-25-26-50/53 S: (1/2-)28-36/37-38-45-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-240-00-0	*cis*-1,2,3,6-tetrahydro-4-methylphthalic anhydride; [1] 1,2,3,6-tetrahydro-4-methylphthalic anhydride; [2] 1,2,3,6-tetrahydro-3-methylphthalic anhydride; [3] tetrahydromethylphthalic anhydride; [4] 1,2,3,6-tetrahydromethylphthalic anhydride; [5] tetrahydro-4-methylphthalic anhydride; [6] 2,3,5,6-tetrahydro-2-methylphthalic anhydride [7]	216-906-6 [1] 222-323-8 [2] 226-247-6 [3] 234-290-7 [4] 247-830-1 [5] 251-823-9 [6] 255-853-3 [7]	1694-82-2 [1] 3425-89-6 [2] 5333-84-6 [3] 11070-44-3 [4] 26590-20-5 [5] 34090-76-1 [6] 42498-58-8 [7]	Xi; R41 R42/43	Xn R: 41-42/43 S: (2-)22-24-26-37/39		C
607-241-00-6	hexahydro-4-methylphthalic anhydride; [1] hexahydromethylphthalic anhydride; [2] hexahydro-1-methylphthalic anhydride; [3] hexahydro-3-methylphthalic anhydride [4]	243-072-0 [1] 247-094-1 [2] 256-356-4 [3] 260-566-1 [4]	19438-60-9 [1] 25550-51-0 [2] 48122-14-1 [3] 57110-29-9 [4]	Xi; R41 R42/43	Xn R: 41-42/43 S: (2-)22-24-26-37/39		C
607-242-00-1	tetrachlorophthalic anhydride	204-171-4	117-08-8	Xi; R41 R42/43 N; R50-53	Xn; N R: 41-42/43-50/53 S: (2-)22-24-26-37/39-60-61		
607-243-00-7	sodium 3,6-dichloro-*o*-anisate; [1] 3,6-dichloro-*o*-anisic acid, compound with 2,2'-iminodiethanol (1:1); [2] 3,6-dichloro-*o*-anisic acid, compound with 2-aminoethanol (1:1) [3]	217-846-3 [1] 246-590-5 [2] 258-527-9 [3]	1982-69-0 [1] 25059-78-3 [2] 53404-28-7 [3]	R52-53	R: 52/53 S: 61		
607-244-00-2	isooctyl acrylate	249-707-8	29590-42-9	Xi; R36/37/38 N; R50-53	Xi; N R: 36/37/38-50/53 S: (2-)26-28-60-61	Xi; R36/37/38: C ≥ 10 %	
607-245-00-8	*tert*-butyl acrylate	216-768-7	1663-39-4	F; R11 Xn; R20/21/22 Xi; R37/38 R43 N; R52-53	F; Xn R: 11-20/21/22-37/38-43-52/53 S: (2-)16-25-37-61		D

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
607-246-00-3	allyl methacrylate; 2-methyl-2-propenoic acid 2-propenyl ester	202-473-0	96-05-9	R10 T; R23 Xn; R21/22 N; R50	T; N R: 10-21/22-23-50 S: (1/2-)36/37-45-61		
607-247-00-9	dodecyl methacrylate	205-570-6	142-90-5	Xi; 36/37/38 N; R50-53	Xi; N R: 36/37/38-50/53 S: (2-)26-28-60-61	Xi; R36/37/38: C ≥ 10 %	
607-248-00-4	naptalam-sodium (ISO); sodium *N*-naphth-1-ylphthalamate	205-073-4	132-67-2	Xn; R22	Xn R: 22 S: (2-)		
607-249-00-X	(1-methyl-1,2-ethanediyl)bis[oxy(methyl-2,1-ethanediyl)] diacrylate	256-032-2	42978-66-5	Xi; R36/37/38 R43 N; R51-53	Xi; N R: 36/37/38-43-51/53 S: (2-)24-37-61	Xi; R36/37/38: C ≥ 10 %	
607-250-00-5	4*H*-3,1-benzoxazine-2,4(1*H*)-dione	204-255-0	118-48-9	Xi; R36 R43	Xi R: 36-43 S: (2-)24-26-37		
607-251-00-0	2-methoxypropyl acetate	274-724-2	70657-70-4	R10 Repr. Cat. 2; R61 Xi; R37	T R: 61-10-37 S: 53-45		
607-252-00-6	lambda-cyhalothrin (ISO); reaction mass of (*S*)-α-cyano-3-phenoxybenzyl(*Z*)-(1*R*)-*cis*-3-(2-chloro-3,3,3-trifluoropropenyl)-2,2-dimethylcyclopropanecarboxylate and (*R*)-α-cyano-3-phenoxybenzyl (*Z*)-(1*S*)-*cis*-3-(2-chloro-3,3,3-trifluoropropenyl)-2,2-dimethylcyclopropanecarboxylate (1:1)	415-130-7	91465-08-6	T+; R26 T; R25 Xn; R21 N; R50-53	T+; N R: 21-25-26-50/53 S: (1/2-)28-36/37/39-38-45-60-61		
607-253-00-1	α-cyano-4-fluoro-3-phenoxybenzyl-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate; cyfluthrin	269-855-7	68359-37-5	T+; R28 T; R23 N; R50-53	T+; N R: 23-28-50/53 S: (1/2-)36/37/39-45-60-61		
607-254-00-7	α-cyano-4-fluoro-3-phenoxybenzyl-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate; beta-cyfluthrin	269-855-7	68359-37-5	T+; R26/28 N; R50-53	T+; N R: 26/28-50/53 S: (1/2-)36/37/39-45-60-61		

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607-255-00-2	fluroxypyr (ISO); 4-amino-3,5-dichloro-6-fluoro-2-pyridyloxyacetic acid	—	69377-81-7	R52-53	R: 52/53 S: 61		
607-256-00-8	azoxystrobin (ISO); methyl (*E*)-2-{2-[6-(2-cyanophenoxy)pyrimidin-4-yloxy]phenyl}-3-methoxyacrylate	—	131860-33-8	T; R23 N; 50-53	T; N R: 23-50/53 S: (1/2-)22-45-60-61		
607-257-00-3	isopropyl propionate	211-300-8	637-78-5	F; R11	F R: 11 S: (2-)16-23-24-29-33		
607-258-00-9	dodecyl 3-(2-(3-benzyl-4-ethoxy-2,5-dioxoimidazolidin-1-yl)-3-(4-methoxybenzoyl)acetamido)-4-chlorobenzoate	403-990-6	70950-45-7	R53	R: 53 S: 61		
607-259-00-4	methyl 2*R*,3*S*-(-)-3-(4-methoxyphenyl)oxiranecarboxylate	404-130-2	105560-93-8	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
607-260-00-X	ethyl 2-(3-nitrobenzylidene)acetoacetate	404-490-0	39562-16-8	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
607-261-00-5	iso(C_{10}-C_{14})alkyl (3,5-di-*tert*-butyl-4-hydroxyphenyl) methylthioacetate	404-800-4	118832-72-7	N; R50-53	N R: 50/53 S: 60-61		
607-262-00-0	7-chloro-1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxoquinoline-3-carboxylic acid	405-050-0	86393-33-1	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
607-263-00-6	potassium iron(III) 1,3-propanediamine-*N*,*N*,*N*',*N*'-tetraacetate hemihydrate	405-680-6	—	E; R2 N; R51-53	E; N R: 2-51/53 S: (2-)35-61		
607-264-00-1	2-chloro-4-(methylsulfonyl)benzoic acid	406-520-8	53250-83-2	Xi; R41	Xi R: 41 S: (2-)26-39		
607-265-00-7	ethyl-2-chloro-2,2-diphenylacetate	406-580-5	52460-86-3	Xi; R38 R52-53	Xi R: 38-52/53 S: (2-)37-61		

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607-266-00-2	reaction mass of: hydroxyaluminium bis[2-hydroxy-3,5-di-*tert*-butylbenzoate]; 3,5-di-*tert*-butyl-salicylic acid	406-890-0	130296-87-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-60-61		
607-267-00-8	*tert*-butyl (5*S*,6*R*,7*R*)-3-bromomethyl-5,8-dioxo-7-(2-(2-phenylacetamido)-5-thia-1-azabicyclo[4.2.0] oct-2-ene-2-carboxylate	407-620-4	33610-13-8	R42/43 R52-53	Xn R: 42/43-52/53 S: (2-)22-24-37-61		
607-268-00-3	2-methylpropyl (*R*)-2-hydroxypropanoate	407-770-0	61597-96-4	Xi; R36	Xi R: 36 S: (2-)26		
607-269-00-9	(*R*)-2-(4-hydroxyphenoxy)propanoic acid	407-960-3	94050-90-5	Xi; R41	Xi R: 41 S: (2-)26-39		
607-270-00-4	3,9-bis(2-(3-(3-*tert*-butyl-4-hydroxy-5-methylphenyl)propionyloxy-1,1-dimethylethyl)-2,4,8,10- tetraoxaspiro[5.5]undecane	410-730-5	90498-90-1	Xn; R21	Xn R: 21 S: (2-)36/37		
607-271-00-X	2-isopropyl-5-methylcyclohexyloxycarbonyloxy-2-hydroxypropane	417-420-9	156324-82-2	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
607-272-00-5	fluroxypyr-meptyl (ISO); methylheptyl, *O*-(4-amino-3,5-dichloro-6-fluoro-2-pyridyloxy) acetate; [1] fluroxypyr-butometyl (ISO); 2-butoxy-1-methylethyl, *O*-(4-amino-3,5-dichloro-6-fluoro-2-pyridyloxy) acetate [2]	279-752-9 [1] - [2]	81406-37-3 [1] 154486-27-8 [2]	N; R50-53	N R: 50/53 S: 60-61		
607-273-00-0	ammonium 7-(2,6-dimethyl-8-(2,2-dimethylbutyryloxy)-1,2,6,7,8,8a-hexahydro-1-naphthyl)-3,5-dihydroxyheptanoate	404-520-2	—	R52-53	R: 52/53 S: 61		
607-274-00-6	2-(*N*-benzyl-*N*-methylamino)ethyl 3-amino-2-butenoate	405-350-1	54527-73-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-275-00-1	sodium benzoyloxybenzene-4-sulfonate	405-450-5	66531-87-1	R43	Xi R: 43 S: (2-)24-37		

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607-276-00-7	bis[(1-methylimidazol)-(2-ethyl-hexanoate)], zinc complex	405-635-0	—	Xi; R38-41 N; R50-53	Xi; N R: 38-41-50/53 S: (2-)26-37/39-60-61		
607-277-00-2	reaction mass of: 2-(hexylthio)ethylamine hydrochloride; sodium propionate	405-720-2	—	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)24-26-37/39-61		
607-278-00-8	reaction mass of isomers of: sodium phenethylnaphthalene-sulfonate; sodium naphthylethylbenzenesulfonate	405-760-0	—	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
607-279-00-3	reaction mass of *n*-octadecylaminodiethyl bis(hydrogen maleate); *n*-octadecylaminodiethyl hydrogen maleate hydrogenphthalate	405-960-8	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-280-00-9	sodium 4-chloro-1-hydroxybutane-1-sulfonate	406-190-5	54322-20-2	Xn; R22 Xi; R36 R43	Xn R: 22-36-43 S: (2-)22-26-36/37		
607-281-00-4	reaction mass of branched and linear C_7-C_9 alkyl 3-[3-(2*H*-benzotriazol-2-yl)-5-(1,1-dimethylethyl)-4-hydroxyphenyl] propionates	407-000-3	127519-17-9	N; R51-53	N R: 51/53 S: 61		
607-282-00-X	2-acetoxymethyl-4-benzyloxybut-1-yl acetate	407-140-5	131266-10-9	R52-53	R: 52/53 S: 61		
607-283-00-5	*E*-ethyl-4-oxo-4-phenylcrotonate	408-040-4	15121-89-8	Xn; R21/22 Xi; R38-41 R43 N; R50-53	Xn; N R: 21/22-38-41-43-50/53 S: (2-)26-36/37/39-60-61		
607-284-00-0	reaction mass of: sodium 3,3'-(1,4-phenylenebis(carbonylimino-3,1-propanediylimino))bis(10-amino-6,13-dichloro-4,11-triphenodioxazinedisulfonate); lithium 3,3'-(1,4-phenylenebis-(carbonylimino-3,1-propanediyl-imino))bis(10-amino-6,13-dichloro)-4,11-triphenodioxazinedisulfonate (9:1)	410-040-4	136213-76-8	N; R51-53	N R: 51/53 S: 61		

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607-285-00-6	reaction mass of: 7-(((3-aminophenyl)sulfonyl)amino)-naphthalene-1,3-disulfonic acid; sodium 7-(((3-aminophenyl)sulfonyl)amino)-naphthalene-1,3-disulfonate; potassium 7-(((3-aminophenyl)sulfonyl)amino)-naphthalene-1,3-disulfonate	410-065-0	—	R43	Xi R: 43 S: (2-)22-24-37		
607-286-00-1	reaction mass of: sodium/potassium 7-[[[3-[[4-((2-hydroxy-naphthyl)azo)phenyl]azo]phenyl]sulfonyl]amino]-naphthalene-1,3-disulfonate	410-070-8	141880-36-6	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
607-287-00-7	*O*'-methyl *O*-(1-methyl-2-methacryloyloxy-ethyl)-1,2,3,6-tetrahydrophthalate	410-140-8	—	R52-53	R: 52/53 S: 61		
607-288-00-2	Tetrasodium (*c*-(3-(1-(3-(*e*-6-dichloro-5-cyanopyrimidin-*f*-yl (methyl)amino)propyl)-1,6-dihydro-2-hydroxy-4-methyl-6-oxo-3-pyridylazo)-4-sulfonatophenylsulfamoyl)phthalocyanine-*a*,*b*,*d*-trisulfonato(6-))nickelato II, where *a* is 1 or 2 or 3 or 4,*b* is 8 or 9 or 10 or 11,*c* is 15 or 16 or 17 or 18, *d* is 22 or 23 or 24 or 25 and where *e* and *f* together are 2 and 4 or 4 and 2 respectively	410-160-7	148732-74-5	Xi; R36 R43 R52-53	Xi R: 36-43-52/53 S: (2-)22-26-36/37-61		
607-289-00-8	3-(3-(4-(2,4-bis(1,1-dimethylpropyl)phenoxy)butylaminocarbonyl-4-hydroxy-1-naphthalenyl)thio)propanoic acid	410-370-9	105488-33-3	R53	R: 53 S: 61		
607-290-00-3	reaction mass (ratio not known) of: ammonium 1-C_{14}-C_{18}-alkyloxycarbonyl-2-(3-allyloxy-2-hydroxypropoxycarbonyl) ethane-1-sulfonate; ammonium 2-C_{14}-C_{18}-alkyloxycarbonyl-1-(3-allyloxy-2-hydroxypropoxycarbonyl)ethane-1-sulfonate	410-540-2	—	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
607-291-00-9	dodecyl-ω-(C_5/C_6-cycloalkyl)alkyl carboxylate	410-630-1	104051-92-5	R53	R: 53 S: 61		
607-292-00-4	reaction mass of: [1-(methoxymethyl)-2-(C_{12}-alkoxy)-ethoxy] acetic acid; [1-(methoxymethyl)-2-(C_{14}-alkoxy)-ethoxy]acetic acid	410-640-6	—	Xi; R38-41 N; R50-53	Xi; N R: 38-41-50/53 S: (2-)26-37/39-60-61		
607-293-00-X	reaction mass of: *N*-aminoethylpiperazonium mono-2,4,6-trimethylnonyldiphenyl ether di-sulfonate; *N*-aminoethylpiperazonium di-2,4,6-trimethylnonyldiphenyl ether di-sulfonate	410-650-0	—	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)26-36/37/39-61		
607-294-00-5	sodium 2-benzoyloxy-1-hydroxyethane-sulfonate	410-680-4	—	R43	Xi R: 43 S: (2-)24-37		

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607-295-00-0	reaction mass of: tetrasodium phosphonoethane-1,2-dicarboxylate; hexasodium phosphonobutane-1,2,3,4-tetracarboxylate	410-800-5	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-296-00-6	reaction mass of: pentaerythriol tetraesters with heptanoic acid and 2-ethylhexanoic acid	410-830-9	—	R53	R: 53 S: 61		
607-297-00-1	(*E-E*)-3,3'-(1,4-phenylenedimethylidene)bis(2-oxobornane-10-sulfonic acid)	410-960-6	92761-26-7	Xi; R41	Xi R: 41 S: (2-)26-39		
607-298-00-7	2-(trimethylammonium)ethoxycarboxybenzene-4-sulfonate	411-010-3	—	R43	Xi R: 43 S: (2-)22-36/37		
607-299-00-2	methyl 3-(acetylthio)-2-methyl-propanoate	411-040-7	97101-46-7	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
607-300-00-6	trisodium [2-(5-chloro-2,6-difluoropyrimidin-4-ylamino)-5-(*b*-sulfamoyl-*c, d*-sulfonatophthalocyanin-*a*-yl-K4,N29,N30,N31,N32-sulfonylamino)benzoato(5-)]cuprate(II) where *a* = 1,2,3,4 *b* = 8,9,10,11 *c* = 15,16,17,18 *d* = 22,23,24,25	411-430-7	—	Xi; R41 R43	Xi R: 41-43 S: (2-)26-36/37/39		
607-301-00-1	reaction mass of: dodecanoic acid; poly(1-7)lactate esters of dodecanoic acid	411-860-5	—	R43 N; R51-53	Xi; N R: 38-41-43-51/53 S: (2-)24-26-37/39-61		
607-302-00-7	reaction mass of: tetradecanoic acid; poly(1-7)lactate esters of tetradecanoic acid	411-910-6	—	Xi; R38-41 R43 N; R51-53	Xi; N R: 38-41-43-51/53 S: (2-)24-26-37/39-61		
607-303-00-2	1-cyclopropyl-6,7-difluoro-1,4-dihydro-4-oxoquinoline-3-carboxylic acid	413-760-7	93107-30-3	Repr. Cat. 3; R62 R52-53	Xn R: 62-52/53 S: (2-)22-36/37-61		
607-304-00-8	fluazifop-butyl (ISO); butyl (*RS*)-2-[4-(5-trifluoromethyl-2-pyridyloxy)phenoxy]propionate	274-125-6	69806-50-4	Repr. Cat. 2; R61 N; R50-53	T; N R: 61-50/53 S: 53-45-60-61		
607-305-00-3	fluazifop-P-butyl (ISO); butyl (*R*)-2-[4-(5-trifluoromethyl-2-pyridyloxy)phenoxy]propionate	—	79241-46-6	Repr. Cat. 3; R63 N; R50-53	Xn; N R: 50/53-63 S: (2-)29-36/37-46-60-61		

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607-306-00-9	chlozolinate (ISO); ethyl (*RS*)-3-(3,5-dichlorophenyl)-5-methyl-2,4-dioxo-oxazolidine-5-carboxylate	282-714-4	84332-86-5	Carc. Cat. 3; R40 N; R51-53	Xn; N R: 40-51/53 S: (2-)36/37-61		
607-307-00-4	vinclozolin (ISO); *N*-3,5-dichlorophenyl-5-methyl-5-vinyl-1,3-oxazolidine-2,4-dione	256-599-6	50471-44-8	Carc. Cat. 3; R40 Repr. Cat. 2; R60-61 R43 N; R51-53	T; N R: 60-61-40-43-51/53 S: 53-45-61		
607-308-00-X	esters of 2,4-D	—	—	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)26-29-36/37-46-60-61		A
607-309-00-5	carfentrazone-ethyl (ISO); ethyl (*RS*)-2-chloro-3-[2-chloro-4-fluoro-5-[4-difluoromethyl-4,5-dihydro-3-methyl-5-oxo-1*H*-1,2,4-triazol-1-yl]phenyl]propionate	—	128639-02-1	N; R50-53	N R: 50/53 S: 60-61		
607-310-00-0	kresoxim-methyl (ISO); methyl (*E*)-2-methoxyimino-[2-(*o*-tolyloxymethyl)phenyl]acetate	—	143390-89-0	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		
607-311-00-6	benazolin-ethyl; ethyl 4-chloro-2-oxo-2*H*-benzothiazole-3-acetate	246-591-0	25059-80-7	N; R51-53	N R: 51/53 S: 61		
607-312-00-1	methoxyacetic acid	210-894-6	625-45-6	Repr. Cat. 2; R60-61 Xn; R22 C; R34	T R: 60-61-22-34 S: 53-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	E
607-313-00-7	neodecanoyl chloride	254-875-0	40292-82-8	T+; R26 Xn; R22 C; R34	T+ R: 22-26-34 S: (1/2-)26-28-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
607-314-00-2	ethofumesate (ISO); (±)-2-ethoxy-2,3-dihydro-3,3-dimethylbenzofuran-5-yl methanesulfonate	247-525-3	26225-79-6	N; R51-53	N R: 51/53 S: 61		

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607-315-00-8	glyphosate (ISO); *N*-(phosphonomethyl)glycine	213-997-4	1071-83-6	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
607-316-00-3	glyphosate-trimesium; glyphosate-trimethylsulfonium	—	81591-81-3	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)36/37-46-61		
607-317-00-9	bis(2-ethylhexyl) phthalate; di-(2-ethylhexyl) phthalate; DEHP	204-211-0	117-81-7	Repr. Cat. 2; R60-61	T R: 60-61 S: 53-45		
607-318-00-4	dibutyl phthalate; DBP	201-557-4	84-74-2	Repr. Cat. 2; R61 Repr. Cat. 3; R62 N; R50	T; N R: 61-50-62 S: 53-45-61		
607-319-00-X	deltamethrin (ISO); (*S*)-α-cyano-3-phenoxybenzyl (1*R*, 3*R*)-3-(2,2-dibromovinyl)-2,2-dimethylcyclopropanecarboxylate	258-256-6	52918-63-5	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)24-28-36/37/39-38-45-60-61		
607-320-00-5	bis[4-(ethenyloxy)butyl] 1,3-benzenedicarboxylate	413-930-0	130066-57-8	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
607-321-00-0	(*S*)-methyl-2-chloropropionate	412-470-8	73246-45-4	R10 Xn; R48/22 Xi; R36	Xn R: 10-36-48/22 S: (2-)23-26-36		
607-322-00-6	4-(4,4-dimethyl-3-oxo-pyrazolidin-1-yl)-benzoic acid	413-120-7	107144-30-9	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-61		
607-323-00-1	2-(1-(2-hydroxy-3,5-di-*tert*-pentyl-phenyl)ethyl)-4,6-di-*tert*-pentylphenyl acrylate	413-850-6	123968-25-2	R53	R: 53 S: 61		
607-324-00-7	reaction mass of: *N*,*N*-di(hydrogenated alkyl C_{14}-C_{18})phtalamic acid; dihydrogenated alkyl (C_{14}-C_{18})amine	413-800-3	—	R53	R: 53 S: 61		

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607-325-00-2	(*S*)-2-chloropropionic acid	411-150-5	29617-66-1	Xn; R21/22 C; R35	C R: 21/22-35 S: (1/2-)23-26-28-36/37/39-45		
607-326-00-8	reaction mass of: isobutyl hydrogen 2-(α-2,4,6-trimethylnon-2-enyl)succinate; isobutyl hydrogen 2-(ß-2,4,6-trimetyhylnon-2-enyl)succinate	410-720-0	141847-13-4	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
607-327-00-3	2-(2-iodoethyl)-1,3-propanediol diacetate	411-780-0	127047-77-2	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)36-61		
607-328-00-9	methyl 4-bromomethyl-3-methoxybenzoate	410-310-1	70264-94-7	Xi; R38-41 R43 N; R50-53	Xi; N R: 38-41-43-50/53 S: (2-)26-36/37/39-60-61		
607-329-00-4	reaction mass of: sodium 2-(C_{12-18}-*n*-alkyl)amino-1,4-butandioate; sodium 2-octadecenyl-amino-1,4-butandioate	411-250-9	—	R43	Xi R: 43 S: (2-)24-26-37/39		
607-330-00-X	(*S*)-2,3-dihydro-1*H*-indole-2-carboxylic acid	410-860-2	79815-20-6	Repr. Cat. 3; R62 Xn; R48/22 R43	Xn R: 43-48/22-62 S: (2-)22-25-26-36/37		
607-331-00-5	reaction mass of: bis(2,2,6,6-tetramethyl-1-octyloxypiperidin-4-yl)-1,10-decanedioate; 1,8-bis[(2,2,6,6-tetramethyl-4-((2,2,6,6-tetramethyl-1-octyloxypiperidin-4-yl)-decan-1,10-dioyl)piperidin-1-yl)oxy] octane	406-750-9	—	R53	R: 53 S: 23-61		
607-332-00-0	cyclopentyl chloroformate	411-460-0	50715-28-1	R10 T; R23 Xn; R22-48/22 Xi; R41 R43	T R: 10-22-23-41-43-48/22 S: (1/2-)26-36/37/39-45		
607-333-00-6	reaction mass of: dodecyl *N*-(2,2,6,6-tetramethylpiperidin-4-yl)-*β*-alaninate; tetradecyl *N*-(2,2,6,6-tetramethylpiperidin-4-yl)-*β*-alaninate	405-670-1	—	Xn; R22-48/22 C; R34 N; R50-53	C; N R: 22-34-48/22-50/53 S: (1/2-)26-36/37/39-45-60-61		

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607-334-00-1	ethyl 1-ethyl-6,7,8-trifluoro-1,4-dihydro-4-oxoquinoline-3-carboxylate	405-880-3	100501-62-0	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
607-335-00-7	methyl (*R*)-2-(4-(3-chloro-5-trifluoromethyl-2-pyridyloxy) phenoxy)propionate	406-250-0	72619-32-0	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
607-336-00-2	4-methyl-8-methylenetricyclo[3.3.1.1^{3,7}]dec-2-yl acetate	406-560-6	122760-85-4	Xi; R38 R43 N; R51-53	Xi; N R: 38-43-51/53 S: (2-)36/37-61		
607-337-00-8	di-*tert*-(C_{12-14})-alkylammonium 2-benzothiazolylthiosuccinate	406-052-4	125078-60-6	R10 Xn; R22 Xi; R38-41 N; R51-53	Xn; N R: 10-22-38-41-51/53 S: (2-)26-37/39-61		
607-338-00-3	2-methylpropyl 2-hydroxy-2-methylbut-3-enoate	406-235-9	72531-53-4	Xi; R36/38	Xi R: 36/38 S: (2-)26-37		
607-339-00-9	2,3,4,5-tetrachlorobenzoylchloride	406-760-3	42221-52-3	Xn; R22 C; R34 R43	C R: 22-34-43 S: (1/2-)26-36/37/39-45		
607-340-00-4	1,3-bis(4-benzoyl-3-hydroxyphenoxy)prop-2-yl acetate	406-990-4	—	N; R51-53	N R: 51/53 S: 61		
607-341-00-X	(9*S*)-9-amino-9-deoxyerythromycin	406-790-7	26116-56-3	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
607-342-00-5	4-chlorobutyl veratrate	410-950-1	69788-75-6	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
607-343-00-0	4,7-methanooctahydro-1*H*-indene-diyldimethyl bis(2-carboxybenzoate)	407-410-2	—	R53	R: 53 S: 61		

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607-344-00-6	reaction mass of: 3-(*N*-(3-dimethylaminopropyl)-(C_{4-8})perfluoroalkylsulfonamido)propionic acid; *N*-[dimethyl-3-(C_{4-8}-perfluoroalkylsulfonamido)propylammonium propionate; 3-(*N*-(3-dimethyl-propylammonium)-(C_{4-8})perfluoroalkylsulfonamido)propionic acid propionate	407-810-7	—	Xn; R48/22	Xn R: 48/22 S: (2-)21-22-36/37		
607-345-00-1	potassium 2-(2,4-dichlorophenoxy)-(*R*)-propionate	413-580-9	113963-87-4	Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43 S: (2-)24-26-37/39		
607-346-00-7	3-icosyl-4-henicosylidene-2-oxetanone	401-210-9	83708-14-9	R53	R: 53 S: 61		
607-347-00-2	sodium (*R*)-2-(2,4-dichlorophenoxy)propionate	413-340-3	119299-10-4	Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43 S: (2-)22-26-36/37/39		
607-348-00-8	magnesium bis((*R*)-2-(2,4-dichlorophenoxy)propionate)	413-360-2	—	Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43 S: (2-)22-26-36/37/39		
607-349-00-3	mono-(tetrapropylammonium) hydrogen 2,2'-dithiobisbenzoate	411-270-8	—	R52-53	R: 52/53 S: 61		
607-350-00-9	bis(4-(1,2-bis(ethoxycarbonyl)ethylamino)-3-methylcyclohexyl)methane	412-060-9	136210-32-7	R43 R52-53	Xi R: 43-52/53 S: (2-)36/37-61		
607-351-00-4	methyl *O*-(4-amino-3,5-dichloro-6-fluoropyridin-2-yloxy) acetate	407-550-4	69184-17-4	N; R51-53	N R: 51/53 S: 20/21-61		
607-352-00-X	4,4'-oxydiphthalic anhydride	412-830-4	1823-59-2	R52-53	R: 52/53 S: 61		
607-353-00-5	reaction mass of: ethyl *exo*-tricyclo[5.2.1.$0^{2,6}$]decane-*endo*-2-carboxylate; ethyl *endo*-tricyclo[5.2.1.02,6]decane-*exo*-2-carboxylate	407-520-0	80657-64-3	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
607-354-00-0	ethyl 2-cyclohexylpropionate	412-280-5	2511-00-4	N; R51-53	N R: 51/53 S: 61		

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607-355-00-6	*p*-tolyl 4-chlorobenzoate	411-530-0	15024-10-9	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
607-356-00-1	ethyl *trans*-2,2,6-trimethylcyclohexanecarboxylate	412-540-8	—	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
607-357-00-7	reaction mass of: *trans*-4-acetoxy-4-methyl-2-propyl-tetrahydro-2*H*-pyran; *cis*-4-acetoxy-4-methyl-2-propyl-tetrahydro-2*H*-pyran	412-450-9	131766-73-9	R43	Xi R: 43 S: (2-)24-37		
607-358-00-2	(1*S*,3*S*,5*R*,6*R*)-(4-nitrophenylmethyl)-1-dioxo-6-phenylacetamido-penam-3-carboxylate	412-670-5	54275-93-3	R42	Xn R: 42 S: (2-)22		
607-359-00-8	(1*S*,4*R*,6*R*,7*R*)-(4-nitrophenylmethyl)3-methylene-1-oxo-7-phenylacetamido-cepham-4-carboxylateido-penam-3-carboxylate	412-800-0	76109-32-5	R42	Xn R: 42 S: (2-)22		
607-360-00-3	sodium 3-acetoacetylamino-4-methoxytolyl-6-sulfonate	411-680-7	133167-77-8	R43	Xi R: 43 S: (2-)24-37		
607-361-00-9	methyl (*R*)-2-(4-hydroxyphenoxy)propionate	411-950-4	96562-58-2	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
607-362-00-4	reaction mass of: (3-methoxy)propylammonium/[tris-(2-hydroxyethyl)]ammonium 2-(2-(bis(2-hydroxyethyl)amino) ethoxycarbonylmethyl)hexadec-4-enoate; (3-methoxy)propylammonium/[tris-(2-hydroxyethyl)]ammonium 2-(2-(bis(2-hydroxyethyl)amino)ethoxycarbonylmethyl) tetradec-4-enoate; (3-methoxy)propylammonium/[tris-(2-hydroxyethyl)]ammonium 2-(3-methoxypropylcarbamoylmethyl)hexadec-4-enoate; (3-methoxy)propylammonium/[tris-(2-hydroxyethyl)]ammonium 2-(3-methoxypropylcarbamoylmethyl)tetradec-4-enoate	413-500-2	—	Xi; R38-41 N; R51-53	Xi; N R: 38-41-51/53 S: (2-)26-37/39-61		
607-363-00-X	methyl-3-methoxyacrylate	412-900-4	5788-17-0	R43	Xi R: 43 S: (2-)24-37		

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607-364-00-5	3-phenyl-7-[4-(tetrahydrofurfuryloxy)phenyl]-1,5-dioxa-s-indacen-2,6-dione	413-330-9	134724-55-3	R53	R: 53 S: 61		
607-365-00-0	2-(2-amino-1,3-thiazol-4-yl)-(*Z*)-2-methoxyiminoacetyl chloride hydrochloride	410-620-7	119154-86-8	Xn; R22 C; R34 R43	C R: 22-34-43 S: (1/2-)22-26-36/37/39-45		
607-366-00-6	3,5-dimethylbenzoyl chloride	413-010-9	6613-44-1	C; R34 R43	C R: 34-43 S: (1/2-)26-36/37/39-45		
607-367-00-1	potassium bis(*N*-carboxymethyl)-*N*-methyl-glycinato-(2-)*N*,*O*,*O*,*N*)-ferrate-(1-) monohydrate	411-640-9	153352-59-1	Xn; R22	Xn R: 22 S: (2-)37		
607-368-00-7	1-(*N*,*N*-dimethylcarbamoyl)-3-*tert*-butyl-5-carbethoxy-methylthio-1*H*-1,2,4-triazole	411-650-3	110895-43-7	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)37-38-45-60-61		
607-369-00-2	reaction mass of: *trans*-(2*R*)-5-acetoxy-1,3-oxathiolane-2-carboxylic acid; *cis*-(2*R*)-5-acetoxy-1,3-oxathiolane-2-carboxylic acid	411-660-8	147027-04-1	Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43 S: (2-)22-24-26-37/39		
607-370-00-8	2-[[2-(acetyloxy)-3-(1,1-dimethyl-ethyl)-5-methylphenyl] methyl]-6-(1,1-dimethylethyl)-4-methylphenol	412-210-3	41620-33-1	N; R50-53	N R: 50/53 S: 60-61		
607-371-00-3	3-ethyl 5-methyl 4-(2-chlorophenyl)-1,4-dihydro-2-[2-(1,3-dihydro-1,3-dioxo-(2*H*)isoindol-2-yl)-ethoxymethyl]-6-methyl-3,5-pyridinedicarboxylate	413-410-3	88150-62-3	R53	R: 53 S: 61		
607-372-00-9	ethoxylated bis phenol A di-(norbornene carboxylate)	412-410-0	—	R52-53	R: 52/53 S: 61		

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607-373-00-4	(±) tetrahydrofurfuryl (*R*)-2-[4-(6-chloroquinoxalin-2-yloxy) phenyloxy]propionate	414-200-4	119738-06-6	Muta. Cat. 3; R68 Repr. Cat. 2; R61 Repr. Cat. 3; R62 Xn; R22-48/22 N; R50-53	T; N R: 61-22-48/22-62-68-50/53 S: 53-45-60-61		E
607-374-00-X	5-amino-2,4,6-triiodo-1,3-benzenedicarbonyldichloride	417-220-1	37441-29-5	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-36/37-61		
607-375-00-5	reaction mass of: *cis*-4-hydroxy-3-(1,2,3,4-tetrahydro-3-(4-(4-trifluoromethylbenzyloxy)phenyl)-1-naphthyl)coumarin; trans-4-hydroxy-3-(1,2,3,4-tetrahydro-3-(4-(4-trifluoro-methylbenzyloxy)phenyl)-1-naphthyl)coumarin	421-960-0	90035-08-8	T+; R26/27/28 T; R48/23/24/25 N; R50-53	T+; N R: 26/27/28-48/23/24/25-50/53 S: (1/2-)28-36/37/39-45-60-61		
607-376-00-0	benzyl 2,4-dibromobutanoate	420-710-8	23085-60-1	Repr. Cat. 3; R62 Xi; R38 R43 N; R50-53	Xn; N R: 38-43-62-50/53 S: (2-)23-36/37-41-60-61		
607-377-00-6	*trans*-4-cyclohexyl-L-proline monohydrochloride	419-160-1	90657-55-9	Repr. Cat. 3; R62 Xn; R22 Xi; R38-41 R43	Xn R: 22-38-41-43-62 S: (2-)22-26-36/37/39		
607-378-00-1	ammonium (*Z*)-α-methoxyimino-2-furylacetate	405-990-1	97148-39-5	F; R11	F R: 11 S: (2-)22-43		
607-379-00-7	reaction mass of: 2-[*N*-(2-hydroxyethyl)stearamido]ethyl stearate; sodium [bis[2-(stearoyloxy)ethyl]amino]methylsulfonate; sodium [bis(2-hydroxyethyl)amino]methylsulfonate; *N*,*N*-bis(2-hydroxyethyl)stearamide	401-230-8		R52-53	R: 52/53 S: 61		

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607-380-00-2	reaction mass of: ammonium-1,2-bis(hexyloxycarbonyl)ethanesulfonate; ammonium-1-hexyloxycarbonyl-2-octyloxycarbonylethanesulfonate; ammonium-2-hexyloxycarbonyl-1-octyloxycarbonylethanesulfonate	407-320-3	—	Xi; R38-41 R52-53	Xi R: 38-41-52/53 S: (2-)26-37/39-61		
607-381-00-8	reaction mass of triesters of 2,2-bis(hydroxymethyl)butanol with C_7-alkanoic acids and 2-ethylhexanoic acid	413-710-4	—	R53	R: 53 S: 61		
607-382-00-3	2-((4-amino-2-nitrophenyl)amino)benzoic acid	411-260-3	117907-43-4	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
607-383-00-9	reaction mass of: 2,2,6,6-tetramethylpiperidin-4-yl-hexadecanoate; 2,2,6,6-tetramethylpiperidin-4-yl-octadecanoate	415-430-8	86403-32-9	Xi; R41 R43 N; R50-53	Xi; N R: 41-43-50/53 S: (2-)24-26-37/39-60-61		
607-384-00-4	reaction mass of: esters of C_{14}-C_{15} branched alcohols with 3,5-di-t-butyl-4-hydroxyphenyl propionic acid; C_{15} branched and linear alkyl 3,5-bis(1,1-dimethylethyl)-4-hydroxybenzenepropanoate; C_{13} branched and linear alkyl 3,5-bis(1,1-dimethylethyl)-4-hydroxybenzenepropanoate	413-750-2	171090-93-0	R53	R: 53 S: 61		
607-385-00-X	Copolymer of vinyl-alcohol and vinyl acetate partially acetilized with 4-(2-(4-formylphenyl)ethenyl)-1-methylpyridinium methylsulfate	414-590-6	125229-74-5	N; R51-53	N R: 51/53 S: 61		
607-386-00-5	reaction mass of: tetradecanoic acid (42.5-47.5 %); poly(1-7)lactate esters of tetradecanoic acid (52.5-57.5 %)	412-580-6	174591-51-6	Xi; R38-41 R43 N; R50-53	Xi; N R: 38-41-43-50/53 S: (2-)24-26-37/39-60-61		
607-387-00-0	reaction mass of: dodecanoic acid (35-40 %); poly(1-7)lactate esters of dodecanoic acid (60-65 %)	412-590-0	58856-63-6	Xi; R38-41 R43 N; R50-53	Xi; N R: 38-41-43-50/53 S: (2-)24-26-37/39-60-61		
607-388-00-6	4-ethylamino-3-nitrobenzoic acid	412-090-2	2788-74-1	Xn; R22 R43 R52-53	Xn R: 22-43-52/53 S: (2-)22-24-37-61		

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607-389-00-1	trisodium *N*,*N*-bis(carboxymethyl)-3-amino-2-hydroxypropionate	414-130-4	119710-96-2	Xn; R22	Xn R: 22 S: (2-)22		
607-390-00-7	1,2,3,4-tetrahydro-6-nitro-quinoxaline	414-270-6	41959-35-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-61		
607-391-00-2	dimethylcyclopropane-1,1-dicarboxylate	414-240-2	6914-71-2	R52-53	R: 52/53 S: 61		
607-392-00-8	2-phenoxyethyl 4-((5-cyano-1,6-dihydro-2-hydroxy-1,4-dimethyl-6-oxo-3-pyridinyl)azo)benzoate	414-260-1	88938-37-8	R53	R: 53 S: 61		
607-393-00-3	3-(*cis*-1-propenyl)-7-amino-8-oxo-5-thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylic acid	415-750-8	106447-44-3	R43	Xi R: 43 S: (2-)22-24-37		
607-394-00-9	5-methylpyrazine-2-carboxylic acid	413-260-9	5521-55-1	Xi; R41	Xi R: 41 S: (2-)26-39		
607-395-00-4	reaction mass of: sodium 1-tridecyl-4-allyl-(2 or 3)-sulfobutanedioate; sodium 1-dodecyl-4-allyl-(2 or 3)-sulfobutanedioate	410-230-7	—	C; R34 R43 N; R51-53	C; N R: 34-43-51/53 S: (1/2-)26-36/37/39-45-61		
607-396-00-X	bis(1,2,2,6,6-pentamethyl-4-piperidinyl) 2-(4-methoxybenzylidene)malonate	414-840-4	147783-69-5	N; R50-53	N R: 50/53 S: 22-60-61		
607-397-00-5	reaction mass of: Ca salicylates (branched C_{10-14} and C_{18-30} alkylated); Ca phenates (branched C_{10-14} and C_{18-30} alkylated); Ca sulfurized phenates (branched C_{10-14} and C_{18-30} alkylated)	415-930-6	—	R43	Xi R: 43 S: (2-)36/37		
607-398-00-0	ethyl *N*-(5-chloro-3-(4-(diethylamino)-2-methylphenylimino)-4-methyl-6-oxo-1,4-cyclohexadienyl)carbamate	414-820-5	125630-94-6	N; R50-53	N R: 50/53 S: 60-61		
607-399-00-6	2,2-dimethyl 3-methyl-3-butenyl propanoate	415-610-6	104468-21-5	Xi; R38 R52-53	Xi R: 38-52/53 S: (2-)37-61		

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607-400-00-X	methyl 3-[[(dibutylamino)thioxomethyl]thio]propanoate	414-400-1	32750-89-3	N; R50-53	N R: 50/53 S: 60-61		
607-401-00-5	ethyl 3-hydroxy-5-oxo-3-cyclohexene-1-carboxylate	414-450-4	88805-65-6	Xi; R38-41 R43	Xi R: 38-41-43 S: (2-)24-26-37/39		
607-402-00-0	methyl *N*-(phenoxycarbonyl)-L-valinate	414-500-5	153441-77-1	R52-53	R: 52/53 S: 61		
607-403-00-6	reaction mass of: bis(1*S*,2*S*,4*S*)-(1-benzyl-4-*tert*-butoxycarboxamido-2-hydroxy-5-phenyl)pentylammonium succinate; isopropyl alcohol	414-810-0	—	Xn; R48/22 Xi; R41 N; R50-53	Xn; N R: 41-48/22-50/53 S: (2-)22-26-36/39-60-61		
607-404-00-1	reaction mass of: ((*Z*)-3,7-dimethyl-2,6-octadienyl)oxycarbonylpropanoic acid; di-((*E*)-3,7-dimethyl-2,6-octadienyl) butandioate; di-((*Z*)-3,7-dimethyl-2,6-octadienyl) butandioate; (*Z*)-3,7-dimethyl-2,6-octadienyl butandioate; ((*E*)-3,7-dimethyl-2,6-octadienyl)oxycarbonylpropanoic acid	415-190-4	—	R43	Xi R: 43 S: (2-)24-37		
607-405-00-7	2-hexyldecyl-*p*-hydroxybenzoate	415-380-7	148348-12-3	N; R51-53	N R: 51/53 S: 61		
607-406-00-2	potassium 2,5-dichlorobenzoate	415-700-5	184637-62-5	Xn; R22 Xi; R41	Xn R: 22-41 S: (2-)26-39		
607-407-00-8	ethyl 2-carboxy-3-(2-thienyl)propionate	415-680-8	143468-96-6	Xi; R38-41 R43	Xi R: 38-41-43 S: (2-)24-26-37/39		
607-408-00-3	potassium *N*-(4-fluorophenyl)glycinate	415-710-1	184637-63-6	Xn; R48/22 Xi; R41 R43 R52-53	Xn R: 41-43-48/22-52/53 S: (2-)22-26-36/37/39-61		
607-409-00-9	reaction mass of: (3*R*)-[1*S*-(1α, 2α, 6β-((2*S*)-2-methyl-1-oxobutoxy)-8aγ)hexahydro-2,6-dimethyl-1-naphthalene]-3,5-dihydroxyheptanoic acid; inert biomass from *Aspergillus terreus*	415-840-7	—	R43 R52-53	Xi R: 43-52/53 S: (2-)36/37-61		

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607-410-00-4	mono[2-(dimethylamino)ethyl]monohydrogen-2-(hexadec-2-enyl)butanedioate and/or mono[2-(dimethylamino)ethyl] monohydrogen-3-(hexadec-2-enyl)butanedioate	415-880-5	779343-34-9	Xi; R38-41 R43 N; R50-53	Xi; N R: 38-41-43-50/53 S: (2-)24-26-37/39-60-61		
607-411-00-X	oxiranemethanol, 4-methylbenzene-sulfonate, (*S*)-	417-210-7	70987-78-9	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xi; R41 R43 N; R51-53	T; N R: 45-41-43-68-51/53 S: 53-45-61		
607-412-00-5	ethyl 2-(1-cyanocyclohexyl)acetate	415-970-4	133481-10-4	Xn; R22-48/22 R52-53	Xn R: 22-48/22-52/53 S: (2-)36/37-61		
607-413-00-0	trans-4-phenyl-L-proline	416-020-1	96314-26-0	Repr. Cat. 3; R62 R43	Xn R: 43-62 S: (2-)22-36/37		
607-414-00-6	tris(2-ethylhexyl)-4,4',4"-(1,3,5-triazine-2,4,6-triyltriimino)tribenzoate	402-070-1	88122-99-0	R53	R: 53 S: 61		
607-415-00-1	poly-(methyl methacrylate)-co-(butylmethacrylate)-co-(4-acryloxybutyl-isopropenyl-α, α-dimethylbenzyl carbamate)-co-(maleicanhydride)	419-590-1	—	F; R11 R43	F; Xi R: 11-43 S: (2-)24-37-43		
607-416-00-7	4-(2-carboxymethylthio)ethoxy-1-hydroxy-5-isobutyloxycarbonylamino-*N*-(3-dodecyloxypropyl)-2-naphthamide	420-730-7	—	N; R50-53	N R: 50/53 S: 60-61		
607-418-00-8	2-ethylhexyl 4-aminobenzoate	420-170-3	26218-04-2	N; R50-53	N R: 50/53 S: 60-61		
607-419-00-3	(3'-carboxymethyl-5-(2-(3-ethyl-3*H*-benzothiazol-2-ylidene)-1-methyl-ethylidene)-4,4'-dioxo-2'-thioxo-(2,5')bithiazolidinyliden-3-yl)-acetic acid	422-240-9	166596-68-5	Xi; R41 R43	Xi R: 41-43 S: (2-)26-36/37/39		
607-420-00-9	2,2-bis(hydroxymethyl)butanoic acid	424-090-1	10097-02-6	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		

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607-421-00-4	cypermethrin *cis*/*trans* +/− 40/60; (*RS*)-α-cyano-3-phenoxybenzyl (1*RS*,3*RS*;1*RS*,3*SR*)-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate	257-842-9	52315-07-8	Xn; R20/22 Xi; R37 N; R50-53	Xn; N R: 20/22-37-50/53 S: (2-)24-36/37/39-60-61		
607-422-00-X	α-cypermethrin	257-842-9	67375-30-8	T; R25 Xn; R48/22 Xi; R37 N; R50-53	T; N R: 25-37-48/22-50/53 S: (2-)36/37/39-45-60-61		
607-423-00-5	esters of mecoprop and of mecoprop-P	—	—	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)13-36/37-60-61		A
607-424-00-0	trifloxystrobin (ISO); (*E*,*E*)-α-methoxyimino-{2-[[[[1-[3-(trifluoromethyl)phenyl] ethylidene]amino]oxy]methyl]benzeneacetic acid methyl ester	—	141517-21-7	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-46-60-61		
607-425-00-6	metalaxyl (ISO); methyl-*N*-(2,6-dimethylphenyl)-*N*-(methoxyacetyl)-DL-alaninate	260-979-7	57837-19-1	Xn; R22 R43 R52-53	Xn R: 22-43-52/53 S: (2-)13-24-37-46-61		
607-426-00-1	1,2-benzenedicarboxylic acid, dipentylester, branched and linear; [1] n-pentyl-isopentylphthalate; [2] di-n-pentyl phthalate; [3] diisopentylphthalate [4]	284-032-2 [1] - [2] 205-017-9 [3] 210-088-4 [4]	84777-06-0 [1] - [2] 131-18-0 [3] 605-50-5 [4]	Repr. Cat. 2; R60-61 N; R50	T; N R: 60-61-50 S: 53-45-61		
607-427-00-7	bromoxynil heptanoate (ISO); 2,6-dibromo-4-cyanophenyl heptanoate	260-300-4	56634-95-8	Repr. Cat. 3; R63 Xn; R20/22 R43 N; R50-53	Xn; N R: 20/22-43-63-50/53 S: (2-)36/37-46-60-61		
607-430-00-3	BBP; benzyl butyl phthalate	201-622-7	85-68-7	Repr. Cat. 2; R61 Repr. Cat. 3; R62 N; R50-53	T; N R: 61-62-50/53 S: 53-45-60-61		

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607-431-00-9	prallethrin (ISO); ETOC; 2-methyl-4-oxo-3-(prop-2-ynyl)cyclopent-2-en-1-yl 2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate	245-387-9	23031-36-9	T; R23 Xn; R22 N; R50-53	T; N R: 22-23-50/53 S: (1/2-)45-60-61		
607-432-00-4	*S*-metolachlor; reaction mass of (*S*)-2-chloro-*N*-(2-ethyl-6-methyl-phenyl)-*N*-(2-methoxy-1-methyl-ethyl)-acetamide (80-100 %); [1] (*R*)-2-chloro-*N*-(2-ethyl-6-methyl-phenyl)-*N*-(2-methoxy-1-methyl-ethyl)-acetamide (0-20 %) [2]	- [1] - [2]	87392-12-9 [1] 178961-20-1 [2]	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
607-433-00-X	cypermethrin *cis*/*trans* +/- 80/20; (*RS*)-α-cyano-3-phenoxybenzyl (1*RS*; 3*RS*; 1*RS*, 3*SR*)-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate	257-842-9	52315-07-8	Xn; R22 Xi; R37/38 R43 N; R50-53	Xn; N R: 22-37/38-43-50/53 S: (2-)36/37/39-60-61		
607-434-00-5	mecoprop-P [1] and its salts; (*R*)-2-(4-chloro-2-methylphenoxy)propionic acid	240-539-0	16484-77-8	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)13-26-37/39-46-61		
607-435-00-0	2*S*-isopropyl-5*R*-methyl-1*R*-cyclohexyl 2,2-dihydroxyacetate	416-810-6	111969-64-3	Xn; R48/22 Xi; R41 N; R51-53	Xn; N R: 41-48/22-51/53 S: (2-)22-26-36/39-61		
607-436-00-6	2-hydroxy-3-(2-ethyl-4-methylimidazoyl)propyl neodecanoate	417-350-9	—	Xi; R38-41 N; R50-53	Xi; N R: 38-41-50/53 S: (2-)26-28-37/39-60-61		
607-437-00-1	3-(4-aminophenyl)-2-cyano-2-propenoic acid	417-480-6	252977-62-1	R43	Xi R: 43 S: (2-)22-24-37		
607-438-00-7	methyl-2-[(aminosulfonyl)methyl]benzoate	419-010-5	112941-26-1	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)22-26		
607-439-00-2	methyl tetrahydro-2-furancarboxylate	420-670-1	37443-42-8	Xi; R41	Xi R: 41 S: (2-)26-39		

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607-440-00-8	methyl 2-aminosulfonyl-6-(trifluoromethyl)pyridine-3-c arboxylate	421-220-7	144740-59-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
607-441-00-3	3-[3-(2-dodecyloxy-5-methylphenylcarbamoyl)-4-hydroxy-1-naphthylthio]propionic acid	421-490-6	167684-63-1	R53	R: 53 S: 57-61		
607-442-00-9	benzyl [hydroxy-(4-phenylbutyl)phosphinyl] acetate	416-050-5	87460-09-1	Xi; R41	Xi R: 41 S: (2-)26-36/39		
607-443-00-4	bis(2,4-di-*tert*-butyl-6-methylphenyl)ethyl phosphate	416-140-4	145650-60-8	R53	R: 53 S: 61		
607-444-00-X	reaction mass of: *cis*-1,4-dimethylcyclohexyl dibenzoate; *trans*-1,4-dimethylcyclohexyl dibenzoate	416-230-3	35541-81-2	R53	R: 53 S: 61		
607-445-00-5	Iron (III) tris(4-methylbenzenesulfonate)	420-960-8	77214-82-5	Xi; R41	Xi R: 41 S: (2-)24-26-39		
607-446-00-0	methyl 2-[4-(2-chloro-4-nitrophenylazo)-3-(1-oxopropyl) amino]phenylaminopropionate	416-240-8	155522-12-6	R43 R53	Xi R: 43-53 S: (2-)22-24-37-61		
607-447-00-6	sodium 4-[4-(4-hydroxyphenylazo)phenylamino]-3-nitroben-zenesulfonate	416-370-5	156738-27-1	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
607-448-00-1	2,3,5,6-tetrafluorobenzoic acid	416-800-1	652-18-6	Xi; R38-41	Xi R: 38-41 S: (2-)22-26-37/39		
607-449-00-7	reaction mass of: 4,4',4"-[(2,4,6-trioxo-1,3,5(2*H*,4*H*,6*H*)-tria-zine-1,3,5-triyl)tris[methylene(3,5,5-trimethyl-3,1-cyclohexa-nediyl)iminocarbonyloxy-2,1-ethanediyl(ethyl)amino]]trisben-zenediazoniumtri[bis(2-methylpropyl)naphthalenesulfonate]; 4,4',4",4"'-[[5,5'-[carbonylbis[imino(1,5,5-trimethyl-3,1-cyclo-hexanediyl)methylene]]-2,4,6-trioxo-1,3,5(2*H*,4*H*,6*H*)-tria-zine-1,1',3,3'-tetrayl]tetrakis[methylene(3,5,5-trimethyl-3,1-cyclohexanediyl)iminocarbonyloxy-2,1-ethanediyl(ethyl) amino]]tetrakisbenzenediazoniumtetra[bis(2-methylpropyl) naphthalenesulfonate]	417-080-1	—	E; R2 R43 N; R50-53	E; Xi; N R: 2-43-50/53 S: (2-)24-37-60-61		

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607-450-00-2	2-mercaptobenzothiazolyl-(Z)-(2-aminothiazol-4-yl)-2-(*tert*-butoxycarbonyl) isopropoxyiminoacetate	419-040-9	89604-92-2	R53	R: 53 S: 61		
607-451-00-8	4-[4-amino-5-hydroxy-3-(4-(2-sulfoxyethylsulfonyl)phenylazo)-2,7-disulfonapht-6-ylazo]-6-[3-(4-amino-5-hydroxy-3-(4-(2-sulfoxyethylsulfonyl)phenylazo)-2,7-disulfonapht-6-ylazo]phenylcarbonylamino]benzenesulfonic acid, sodium salt	417-640-5	161935-19-9	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
607-453-00-9	4-benzyl-2,6-dihydroxy-4-aza-heptylene bis(2,2-dimethyloctanoate)	418-100-1	172964-15-7	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
607-454-00-4	reaction mass of: *trans*-2-(1-methylethyl)-1,3-dioxane-5-carboxylic acid; *cis*-2-(1-methylethyl)-1,3-dioxane-5-carboxylic acid	418-170-3	116193-72-7	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)25-26-39-61		
607-455-00-X	1-amino-4-(3-[4-chloro-6-(2,5-di-sulfophenylamino)-1,3,5-triazin-2-ylamino]-2,2-dimethyl-propylamino)-anthraquinone-2-sulfonic acid, sodium/lithiumsalt	419-520-8	172890-93-6	R43	Xi R: 43 S: (2-)22-24-37		
607-456-00-5	3-amino-4-chlorobenzoic acid, hexadecyl ester	419-700-6	143269-74-3	N; R51-53	N R: 51/53 S: 61		
607-457-00-0	tetrasodium dihydrogen 1,1''-dihydroxy-8,8''-[p-phenylbis(imino-{6-[4-(2-aminoethyl)piperazin-1-yl]}-1,3,5-triazine-4,2-diyl-imino)]bis(2,2'-azonaphthalene-1',3,6-trisulfonate)	420-350-1	172277-97-3	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
607-458-00-6	reaction mass of: 2-ethyl-[2,6-dibromo-4-[1-[3,5-dibromo-4-(2-hydroxyethoxy)phenyl]-1-methylethyl]phenoxy]propenoate; 2,2'-diethyl-[4,4'-bis(2,6-dibromophenoxy)-1-methylethylidene] dipropenoate; 2,2'-[(1-methylethylidene)bis[[2,6-dibromo-4,1-phenylene)oxy]ethanol]]	420-850-1	—	N; R51-53	N R: 51/53 S: 61		
607-459-00-1	isopentyl 4-{2-[5-cyano-1,2,3,6-tetrahydro-1-(2-isopropoxyethoxy-carbonylmethyl)-4-methyl-2,6-dioxo-3-pyridylidene]hydrazino}benzoate	418-930-4	—	R53	R: 53 S: 61		
607-460-00-7	3-tridecyloxy-propyl-ammonium 9-octadecenoate	418-990-1	778577-53-0	Xn; R48/22 Xi; R36/38 N; R50-53	Xn; N R: 36/38-48/22-50/53 S: (2-)23-26-37/39-60-61		

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607-461-00-2	reaction mass of: pentasodium 2-{4-{3-methyl-4-[6-sulfonato-4-(2-sulfonato-phenylazo)-naphthalen-1-ylazo]-phenylamino}-6-[3-(2-sulfato-ethanesulfonyl)-phenylamino]-1,3,5-triazin-2-ylamino}-benzene-1,4-disulfonate; pentasodium 2-{4-{3-methyl-4-[7-sulfonato-4-(2-sulfonato-phenylazo)-naphthalen-1-ylazo]-phenylamino}-6-[3-(2-sulfato-ethanesulfonyl)-phenylamino]-1,3,5-triazin-2-ylamino}-benzene-1,4-disulfonate	421-160-1	—	R52-53	R: 52/53 S: 61		
607-462-00-8	reaction mass of: 1-hexyl acetate; 2-methyl-1-pentyl acetate; 3-methyl-1-pentyl acetate; 4-methyl-1-pentyl acetate; other mixed linear and branched C_6-alkyl acetates	421-230-1	88230-35-7	N; R51-53	N R: 51/53 S: 61		
607-463-00-3	3-(phenothiazin-10-yl)propionic acid	421-260-5	362-03-8	N; R51-53	N R: 51/53 S: 24/25-61		
607-464-00-9	reaction mass of: 7-chloro-1-ethyl-6-fluoro-1,4-dihydro-4-oxo-quinoline-3-carboxylic acid; 5-chloro-1-ethyl-6-fluoro-1,4-dihydro-4-oxo-quinoline-3-carboxylic acid	421-280-4		R52-53	R: 52/53 S: 61		
607-465-00-4	tris(2-hydroxyethyl)ammonium 7-{4-[4-(2-cyanoamino-4-hydroxy-6-oxidopyrimidin-5-ylazo)benzamido]-2-ethoxy-phenylazo}naphthalene-1,3-disulfonate	421-440-3	778583-04-3	R52-53	R: 52/53 S: 61		
607-466-00-X	reaction mass of: phenyl 1-(1-[2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl]-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate; phenyl 2-(1-(2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl)-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate; phenyl 3-(1-(2-chloro-5-(hexadecyloxycarbonyl)phenylcarbamoyl)-3,3-dimethyl-2-oxobutyl)-1*H*-2,3,3a,7a-tetrahydrobenzotriazole-5-carboxylate	421-480-1	—	N; R51-53	N R: 51/53 S: 37/39-61		
607-467-00-5	1,1,3,3-tetrabutyl-1,3-ditinoxydicaprylate	419-430-9	56533-00-7	Xn; R21/22-48/22 C; R34 N; R50-53	C; N R: 21/22-34-48/22-50/53 S: (1/2-)26-36/37/39-45-60-61		

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607-468-00-0	reaction mass of: monosodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; disodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; trisodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate; tetrasodium 4-((4-(5-sulfonato-2-methoxyphenylamino)-6-chloro-1,3,5-triazine-2-yl)amino)-2-((1,4-dimethyl-6-oxido-2-oxo-5-sulfonatomethyl-1,2-dihydropyridine-3-yl)azo)benzenesulfonate	419-450-8	—	R43	Xi R: 43 S: (2-)22-24-37		
607-469-00-6	disodium 7-((4,6-bis(3-diethylaminopropylamino)-1,3,5-triazine-2-yl)amino)-4-hydroxy-3-(4-(4-sulfonatophenylazo)phenylazo)-2-naphthalene sulfonate	419-460-2	120029-06-3	R52-53	R: 52/53 S: 61		
607-470-00-1	potassium sodium 6,13-dichloro-3,10-bis{2-[4-[3-(2-hydroxysulphonyloxyethanesulfonyl)phenylamino]-6-(2,5-disulfonatophenylamino)-1,3,5-triazin-2-ylamino]ethylamino}benzo[5,6][1,4]oxazino[2,3-b]phenoxazine-4,11-disulfonate	414-100-0	154336-20-6	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)39-22-26-61		
607-472-00-2	ammonium iron(III) trimethylenediaminetetraacetate hemihydrate	400-660-3	111687-36-6	N; R51-53	N R: 51/53 S: 61		
607-474-00-3	(4-(4-(4-dimethylaminobenzyliden-1-yl)-3-methyl-5-oxo-2-pyrazolin-1-yl)benzoic acid	410-430-4	117573-89-4	R53	R: 53 S: 61		
607-475-00-9	reaction mass of: tetrasodium 7-(4-[4-chloro-6-[methyl-(3-sulfonatophenyl)amino]-1,3,5-triazin-2-ylamino]-2-ureidophenylazo)naphthalene-1,3,6-trisulfonate; tetrasodium 7-(4-[4-chloro-6-[methyl-(4-sulfonatophenyl)amino]-1,3,5-triazin-2-ylamino]-2-ureidophenylazo)naphthalene-1,3,6-trisulfonate (1:1)	412-940-2	148878-18-6	R43	Xi R: 43 S: (2-)22-24-37		
607-476-00-4	trisodium *N*,*N*-bis(carboxymethyl)-β-alanine	414-070-9	129050-62-0	C; R34 R52-53	C R: 34-52/53 S: (1/2-)26-36/37/39-45-61		
607-478-00-5	tetramethylammonium hydrogen phthalate	416-900-5	79723-02-7	T; R25 Xn; R48/22 N; R50	T; N R: 25-48/22-50 S: (1/2-)25-36-45-61		

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607-479-00-0	hexadecyl 4-chloro-3-[2-(5,5-dimethyl-2,4-dioxo-1,3-oxazolidin-3-yl)-4,4-dimethyl-3-oxopentamido]benzoate	418-550-9	168689-49-4	R53	R: 53 S: 61		
607-480-00-6	1,2-benzenedicarboxylic acid; di-C_{7-11}-branched and linear alkylesters	271-084-6	68515-42-4	Repr. Cat. 2; R61 Repr. Cat. 3; R62	T R: 61-62 S: 53-45		
607-487-00-4	reaction mass of: disodium 4-(3-ethoxycarbonyl-4-(5-(3-ethoxycarbonyl-5-hydroxy-1-(4-sulfonatophenyl)pyrazol-4-yl)penta-2,4-dienylidene)-4,5-dihydro-5-oxopyrazol-1-yl)benzenesulfonate; trisodium 4-(3-ethoxycarbonyl-4-(5-(3-ethoxycarbonyl-5-oxido-1-(4-sulfonatophenyl)pyrazol-4-yl)penta-2,4-dienylidene)-4,5-dihydro-5-oxopyrazol-1-yl)benzenesulfonate	402-660-9	—	Repr. Cat. 2; R61 R52-53	T R: 61-52/53 S: 53-45-61		
607-488-00-X	ethyl (2-acetylamino-5-fluoro-4-isothiocyanatophenoxy)acetate	414-210-9	147379-38-2	N; R50-53	N R: 50/53 S: 60-61		
607-489-00-5	reaction mass of: 2-ethylhexyl linolenate, linoleate and oleate; 2-ethylhexyl epoxyoleate; 2-ethylhexyl diepoxylinoleate; 2-ethylhexyl triepoxylinolenate	414-890-7	71302-79-9	R43	Xi R: 43 S: (2-)24-37		
607-490-00-0	*N*-[2-hydroxy-3-(C_{12-16}-alkyloxy)propyl]-*N*-methyl glycinate	415-060-7	—	Xi; R41 R43	Xi R: 41-43 S: (2-)24-26-37/39		
607-492-00-1	2-(1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy)-2-methyl propyl propanoate	415-490-5	141773-73-1	N; R51-53	N R: 51/53 S: 61		
607-493-00-7	methyl (3a*R*,4*R*,7a*R*)-2-methyl-4-(1*S*,2*R*,3-triacetoxypropyl)-3a,7a-dihydro-4*H*-pyrano[3,4-d]oxazole-6-carboxylate	415-670-3	78850-37-0	Xi; R41	Xi R: 41 S: (2-)26-39		
607-494-00-2	bis(2-ethylhexyl)octylphosphonate	417-170-0	52894-02-7	N; R50-53	N R: 50/53 S: 60-61		
607-495-00-8	sodium 4-sulfophenyl-6-((1-oxononyl)amino)hexanoate	417-550-6	168151-92-6	R43	Xi R: 43 S: (2-)24-37		

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607-496-00-3	2,2'-methylenebis(4,6-di-*tert*-butyl-phenyl)-2-ethylhexyl phosphite	418-310-3	126050-54-2	R53	R: 53 S: 61		
607-497-00-9	cerium oxide isostearate	419-760-3	—	R53	R: 53 S: 61		
607-498-00-4	(*E*)-3,7-dimethyl-2,6-octadienylhexadecanoate	421-370-3	3681-73-0	Xi; R38 R53	Xi R: 38-53 S: (2-)37-61		
607-499-00-X	bis(dimethyl-(2-hydroxyethyl)ammonium) 1,2-ethanediyl-bis (2-hexadecenylsuccinate)	421-660-1	—	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24-26-37/39-61		
607-500-00-3	calcium 2,2,bis[(5-tetrapropylene-2-hydroxy)phenyl]ethanoate	421-670-4	—	Xi; R38 N; R50-53	Xi; N R: 38-50/53 S: (2-)37-60-61		
607-501-00-9	reaction mass of: triphenylthiophosphate and tertiary butylated phenyl derivatives	421-820-9	192268-65-8	R53	R: 53 S: 61		
607-502-00-4	(*N*-benzyl-*N*,*N*,*N*-tributyl)ammonium 4-dodecylbenzenesulfonate	422-200-0	178277-55-9	C; R34 Xn; R22 N; R51-53	C; N R: 22-34-51/53 S: (1/2-)26-36/37/39-45-61		
607-503-00-X	2,4,6-tri-*n*-propyl-2,4,6-trioxo-1,3,5,2,4,6-trioxatriphosphorinane	422-210-5	68957-94-8	C; R34	C R: 34 S: (1/2-)26-36/37/39-45		
607-505-00-0	pentasodium 7-(4-(4-(5-amino-4-sulfonato-2-(4-((2-(sulfonato-ethoxy)sulfonyl)phenylazo)phenylamino)-6-chloro-1,3,5-triazin-2-yl)amino-2-ureidophenylazo)naphtalene-1,3,6-trisulfonate	422-930-1		R52-53	R: 52/53 S: 22-61		
607-506-00-6	reaction mass of: strontium (4-chloro-2-((4,5-dihydro-3-methyl-5-oxo-1-(3-sulfonatophenyl)-1*H*-pyrazol-4-yl)azo)-5-methyl)benzenesulfonate; disodium (4-chloro-2-((4,5-dihydro-3-methyl-5-oxo-1-(3-sulfonatophenyl)-1*H*-pyrazol-4-yl)azo)-5-methyl)benzenesulfonate	422-970-8		N; R51-53	N R: 51/53 S: 22-61		

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607-507-00-1	potassium,sodium 2,4-diamino-3-[4-(2-sulfonatoethoxysulfonyl)phenylazo]-5-[4-(2-sulfonatoethoxysulfonyl)-2-sulfonatophenylazo]-benzenesulfonate	422-980-2	187026-95-5	Xi; R41	Xi R: 41 S: (2-)22-26-39		
607-508-00-7	disodium 3,3'-[iminobis[sulfonyl-4,1-phenylene-(5-hydroxy-3-methylpyrazole-1,4-diyl)azo-4,1-phenylenesulfonylimino-(4-amino-6-hydroxypyrimidine-2,5-diyl)azo-4,1-phenylene-sulfonylimino(4-amino-6-hydroxypyrimidine-2,5-diyl)azo]bis (benzenesulfonate)]	423-110-4	—	Xi; R41	Xi R: 41 S: (2-)22-26-39		
607-512-00-9	trisodium 2,4-diamino-3,5-bis-[4-(2-sulfonatoethoxy)sulfonyl)phenylazo]benzenesulfonate	423-970-0	182926-43-8	R52-53	R: 52/53 S: 22-61		
607-513-00-4	reaction mass of: Trisodium 4-benzoylamino-6-(6-ethenesulfonyl-1-sulfato-naphthalen-2-ylazo)-5-hydroxynaphthalene-2,7-disulfonate; 5-(benzoylamino)-4-hydroxy-3-((1-sulfo-6-((2-(sulfooxy)ethyl) sulfonyl)-2-naphthyl)azo)naphthalene-2,7-disulfonic acid sodium salt; 5-(benzoylamino)-4-hydroxy-3-((1-sulfo-6-((2-(sulfooxy)ethyl) sulfonyl)-2-naphthyl)azo)naphthalene-2,7-disulfonic acid	423-200-3	—	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)22-26-36/37/ 39-61		
607-515-00-5	reaction mass of: disodium hexyldiphenyl ether disulphonate; disodium dihexyldiphenyl ether disulphonate	429-650-7	147732-60-3	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
607-516-00-0	*N*,*N*'-bis(trifluoroacetyl)-*S*,*S*'-bis L-homocysteine	429-670-6	105996-54-1	Xi; R41 R43	Xi R: 41-43 S: (2-)24-26-37/39		
607-517-00-6	(*S*)-α-(acetylthio)benzenepropanoic acid	430-300-0	76932-17-7	Xn; R22 Xi; R41 R43	Xn R: 22-41-43 S: (2-)22-26-36/37/39		
607-526-00-5	cartap (ISO); 1,3-bis(carbamoylthio)-2-(dimethylamino)propane	—	15263-53-3	N; R50-53	N R: 50/53 S: 60-61		

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607-527-00-0	reaction mass of: 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-tridecafluorooctyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heptdecafluorodecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heneicosafluorododecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-tridecafluorooctyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-pentacosafluorotetradecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-heptadecafluorodecyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heptadecafluorodecyl)dodecanedioate; 1-(1'*H*,1'*H*,2'*H*,2'*H*-heptadecafluorodecyl)-12-(1"*H*,1"*H*,2"*H*,2"*H*-heneicosafluorododecyl)dodecanedioate	423-180-6	—	Xn; R48/22	Xn R: 48/22 S: (2-)36		
608-001-00-3	acetonitrile; cyanomethane	200-835-2	75-05-8	F; R11 Xn; R20/21/22 Xi; R36	F; Xn R: 11-20/21/22-36 S: (2-)16-36/37		
608-002-00-9	trichloroacetonitrile	208-885-7	545-06-2	T; R23/24/25 N; R51-53	T; N R: 23/24/25-51/53 S: (1/2-)45-61		
608-003-00-4	acrylonitrile	203-466-5	107-13-1	F; R11 Carc. Cat. 2; R45 T; R23/24/25 Xi; R37/38-41 R43 N; R51-53	F; T; N R: 45-11-23/24/25-37/38-41-43-51/53 S: - 53-45-61	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,2 % ≤ C < 1 %	D E
608-004-00-X	2-hydroxy-2-methylpropionitrile; 2-cyanopropan-2-ol; acetone cyanohydrin	200-909-4	75-86-5	T+; R26/27/28 N; R50-53	T+; N R: 26/27/28-50/53 S: (1/2-)7/9-27-45-60-61		
608-005-00-5	*n*-butyronitrile	203-700-6	109-74-0	R10 ⊗ T; R23/24/25	T R: 10-23/24/25 S: (1/2-)45		

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608-006-00-0	bromoxynil (ISO) 3,5-dibromo-4-hydroxybenzonitrile; bromoxynil phenol	216-882-7	1689-84-5	Repr. Cat. 3; R63 T+; R26 T; R25 R43 N; R50-53	T+; N R: 25-26-43-63-50/53 S: (1/2-)27/28-36/37-45-63-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
608-007-00-6	ioxynil (ISO) 4-hydroxy-3,5-diiodobenzonitrile	216-881-1	1689-83-4	Repr. Cat. 3; R63 T; R23/25 Xn; R21-48/22 Xi; R36 N; R50-53	T; N R: 21-23/25-36-48/22-63-50/53 S: (1/2-)36/37-45-60-61-63	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
608-008-00-1	chloroacetonitrile	203-467-0	107-14-2	T; R23/24/25 N; R51-53	T; N R: 23/24/25-51/53 S: (1/2-)45-61		
608-009-00-7	malononitrile	203-703-2	109-77-3	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)23-27-45-60-61		
608-010-00-2	methacrylonitrile; 2-methyl-2-propene nitrile	204-817-5	126-98-7	F; R11 T; R23/24/25 R43	F; T R: 11-23/24/25-43 S: (1/2-)9-16-18-29-45	T; R23/24/25: C ≥ 1 % Xn; R20/21/22: 0,2 % ≤ C < 1 % R43: C ≥ 0,2 %	D
608-011-00-8	oxalonitrile; cyanogen	207-306-5	460-19-5	F; R11 ⊗ T; R23 N; R50-53	F; T; N R: 11-23-50/53 S: (1/2-)23-45-60-61		
608-012-00-3	benzonitrile	202-855-7	100-47-0	Xn; R21/22	Xn R: 21/22 S: (2-)23		
608-013-00-9	2-chlorobenzonitrile	212-836-5	873-32-5	Xn; R21/22 Xi; R36	Xn R: 21/22-36 S: (2-)23		

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608-014-00-4	chlorothalonil (ISO); tetrachloroisophthalonitrile	217-588-1	1897-45-6	Carc. Cat. 3; R40 T+; R26 Xi; R41 Xi; R37 R43 N; R50-53	T+; N R: 26-37-40-41-43-50/53 S: (2-)28-36/37/39-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
608-015-00-X	dichlobenil (ISO); 2,6-dichlorobenzonitrile	214-787-5	1194-65-6	Xn; R21 N; R51-53	Xn; N R: 21-51/53 S: (2-)36/37-61		
608-016-00-5	1,4-Dicyano-2,3,5,6-tetra-chloro-benzene	401-550-8	1897-41-2	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
608-017-00-0	bromoxynil octanoate (ISO); 2,6-dibromo-4-cyanophenyl octanoate	216-885-3	1689-99-2	Repr. Cat. 3; R63 T; R23 Xn; R22 R43 N; R50-53	T; N R: 22-23-43-63-50/53 S: (1/2-)36/37-45-63-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
608-018-00-6	ioxynil octanoate (ISO); 4-cyano-2,6-diiodophenyl octanoate	223-375-4	3861-47-0	Repr. Cat. 3; R63 T; R25 Xi; R36 R43 N; R50-53	T; N R: 25-36-43-63-50/53 S: (1/2-)26-36/37-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
608-019-00-1	2,2'-dimethyl-2,2'-azodipropiononitrile; ADZN	201-132-3	78-67-1	E; R2 F; R11 Xn; R20/22 R52-53	E; Xn R: 2-11-20/22-52/53 S: (2-)39-41-47-61		
608-021-00-2	3-(2-(diaminomethyleneamino)thiazol-4-ylmethylthio)propionitrile	403-710-2	76823-93-3	Xn; R22 R43	Xn R: 22-43 S: (2-)22-24-37		
608-022-00-8	3,7-dimethyloctanenitrile	403-620-3	40188-41-8	Xi; R38 R43 N; R51-53	Xi; N R: 38-43-51/53 S: (2-)36/37-61		

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608-023-00-3	fenbuconazole (ISO) 4-(4-chlorophenyl)-2-phenyl-2-[(1*H*-1,2,4-triazol-1-yl) methyl]butanenitrile	406-140-2	114369-43-6	N; R50-53	N R: 50/53 S: 60-61		
608-024-00-9	2-(4-(*N*-butyl-*N*-phenethylamino)phenyl)ethylene-1,1,2-tricarbonitrile	407-650-8	97460-76-9	R53	R: 53 S: 61		
608-025-00-4	2-nitro-4,5-bis(benzyloxy)phenylacetonitrile	410-970-0	117568-27-1	R53	R: 53 S: 61		
608-026-00-X	3-cyano-3,5,5-trimethylcyclohexanone	411-490-4	7027-11-4	Xn; R22-48/22 R43 R52-53	Xn R: 22-43-48/22-52/53 S: (2-)36/37-61		
608-027-00-5	reaction mass of: 3-(4-ethylphenyl)-2,2-dimethylpropanenitrile; 3-(2-ethylphenyl)-2,2-dimethylpropanenitrile; 3-(3-ethylphenyl)-2,2-dimethylpropanenitrile	412-660-0	—	N; R51-53	N R: 51/53 S: 61		
608-028-00-0	4-(2-cyano-3-phenylamino-acryloyloxymethyl)-cyclohexylmethyl 2-cyano-3-phenylamino)-acrylate	413-510-7	147374-67-2	Xn; R48/20/21 R43 N; R51-53	Xn; N R: 43-48/20/21-51/53 S: (2-)36/37-61		
608-029-00-6	1,2-dihydro-6-hydroxy-4-methyl-1-[3-(1-methylethoxy)propyl]-2-oxo-3-pyridinecarbonitrile	411-990-2	68612-94-2	R43	Xi R: 43 S: (2-)24-37		
608-030-00-1	*N*-acetyl-*N*-[5-cyano-3-(2-dibutylamino-4-phenylthyazol-5-yl-methylene)-4-methyl-2,6-dioxo-1,2,3,6-tetrahydropyridin-1-yl]benzamide	412-340-0	147741-93-3	N; R50-53	N R: 50/53 S: 60-61		
608-031-00-7	2-benzyl-2-methyl-3-butenitrile	407-870-4	97384-48-0	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
608-033-00-8	*N*-butyl-3-(2-chloro-4-nitrophenylhydrazono)-1-cyano-2-methylprop-1-ene-1,3-dicarboximide	407-970-8	75511-91-0	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		

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608-034-00-3	chlorfenapyr; 4-bromo-2-(4-chlorophenyl)-1-ethoxymethyl-5-trifluoro-methylpyrrole-3-carbonitrile	—	122453-73-0	T; R23 Xn; R22 N; R50-53	T; N R: 22-23-50/53 S: (1/2-)13-36/37-45-60-61		
608-035-00-9	(±)-α-[(2-acetyl-5-methylphenyl)-amino]-2,6-dichloroben-zene-aceto-nitrile	419-290-9	—	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
608-036-00-4	3-(2-{4-[2-(4-cyanophenyl)vinyl]phenyl}vinyl)benzonitrile	419-060-8	79026-02-1	R53	R: 53 S: 61		
608-037-00-X	reaction mass of: (*E*)-2,12-tridecadiennitrile; (*E*)-3,12-tridecadiennitrile; (*Z*)-3,12-tridecadiennitrile	422-190-8		N; R50-53	N R: 50/53 S: 60-61		
608-038-00-5	2,2,4-trimethyl-4-phenyl-butane-nitrile	422-580-8	75490-39-0	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
608-039-00-0	2-phenylhexanenitrile	423-460-8	3508-98-3	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)23-60-61		
608-040-00-6	4,4'-dithiobis(5-amino-1-(2,6-dichloro-4-(trifluoromethyl) phenyl)-1*H*-pyrazole-3-carbonitrile)	423-490-1	130755-46-3	N; R50-53	N R: 50/53 S: 60-61		
608-041-00-1	4'-((2-butyl-4-oxo-1,3-diazaspiro[4.4]non-1-ene-3-yl)methyl) (1,1'-biphenyl)-2-carbonitrile	423-500-4	138401-24-8	N; R50-53	N R: 50/53 S: 60-61		
608-043-00-2	3-(*cis*-3-hexenyloxy)propanenitril	415-220-6	142653-61-0	T; R23 Xn; R22 N; R50-53	T; N R: 22-23-50/53 S: (1/2-)13-36/37-45-60-61		
608-065-00-2	salts of bromoxynil with the exception of those specified elsewhere in this Annex	—	—	Repr. Cat. 3; R63 T+; R26 T; R25 R43 N; R50-53	T+; N R: 25-26-43-50/53 S: (1/2-)27/28-36/37-45-63-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	A

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608-066-00-8	salts of ioxynil with the exception of those specified elsewhere in this Annex	—	—	Repr. Cat. 3; R63 T; R23/25 Xn; R21-48/22 Xi; R36 N; R50-53	T; N R: 21-23/25-36-48/22-63-50/53 S: (1/2-)36/37-45-63-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	A
609-001-00-6	1-nitropropane	203-544-9	108-03-2	R10 Xn; R20/21/22	Xn R: 10-20/21/22 S: (2-)9	Xn; R20/21/22: C ≥ 5 %	
609-002-00-1	2-nitropropane	201-209-1	79-46-9	R10 Carc. Cat. 2; R45 Xn; R20/22	T R: 45-10-20/22 S: 53-45		E
609-003-00-7	nitrobenzene	202-716-0	98-95-3	Carc. Cat. 3; R40 Repr. Cat. 3; R62 T; R23/24/25-48/23/24 N; R51-53	T; N R: 23/24/25-40-48/23/24-51/53-62 S: (1/2-)28-36/37-45-61		
609-004-00-2	dinitrobenzene; [1] 1,4-dinitrobenzene; [2] 1,3-dinitrobenzene; [3] 1,2-dinitrobenzene [4]	246-673-6 [1] 202-833-7 [2] 202-776-8 [3] 208-431-8 [4]	25154-54-5 [1] 100-25-4 [2] 99-65-0 [3] 528-29-0 [4]	T+; R26/27/28 R33 N; R50-53	T+; N R: 26/27/28-33-50/53 S: (1/2-)28-36/37-45-60-61		
609-005-00-8	1,3,5-trinitrobenzene	202-752-7	99-35-4	E; R2 ⊗ T+; R26/27/28 R33 N; R50-53	E; T+; N R: 2-26/27/28-33-50/53 S: (1/2-)35-45-60-61		
609-006-00-3	4-nitrotoluene	202-808-0	99-99-0	T; R23/24/25 R33 N; R51-53	T; N R: 23/24/25-33-51/53 S: (1/2-)28-37-45-61		

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609-007-00-9	2,4-dinitrotoluene; dinitrotoluene, technical grade; [1] dinitrotoluene [2]	204-450-0 [1] 246-836-1 [2]	121-14-2 [1] 25321-14-6 [2]	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 N; R51-53	T; N R: 45-23/24/25-48/22-62-68-51/53 S: 53-45-61		E
609-008-00-4	2,4,6-trinitrotoluene; TNT	204-289-6	118-96-7	E; R2 T; R23/24/25 R33 N; R51-53	E; T; N R: 2-23/24/25-33-51/53 S: (1/2-)35-45-61		
609-009-00-X	2,4,6-trinitrophenol; picric acid	201-865-9	88-89-1	E; R2 ⊗ R4 T; R23/24/25	E; T R: 2-4-23/24/25 S: (1/2-)28-35-37-45		
609-010-00-5	salts of picric acid	—	—	E; R3 T; R23/24/25	E; T R: 3-23/24/25 S: (1/2-)28-35-37-45		A
609-011-00-0	2,4,6-trinitroanisole	—	606-35-9	E; R2 Xn; R20/21/22 N; R51-53	E; Xn; N R: 2-20/21/22-51/53 S: (2-)35-61		
609-012-00-6	2,4,6-trinitro-*m*-cresol	210-027-1	602-99-3	E; R2 R4 Xn; R20/21/22	E; Xn R: 2-4-20/21/22 S: (2-)35		
609-013-00-1	2,4,6-trinitro-*m*-xylene	211-187-5	632-92-8	E; R2 Xn; R20/21/22 R33	E; Xn R: 2-20/21/22-33 S: (2-)35		
609-015-00-2	4-nitrophenol; *p*-nitrophenol	202-811-7	100-02-7	Xn; R20/21/22 R33	Xn R: 20/21/22-33 S: (2-)28		

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609-016-00-8	dinitrophenol (reaction mass of isomers); [1] 2,4(or 2,6)-dinitrophenol [2]	247-096-2 [1] 275-732-9 [2]	25550-58-7 [1] 71629-74-8 [2]	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)28-37-45-60-61		
609-018-00-9	2,4,6-trinitroresorcinol; styphnic acid	201-436-6	82-71-3	E; R2 ⊗ R4 Xn; R20/21/22	E; Xn R: 2-4-20/21/22 S: (2-)35		
609-019-00-4	lead 2,4,6-trinitro-*m*-phenylene dioxide; lead 2,4,6-trinitroresorcinoxide; lead styphnate	239-290-0	15245-44-0	E; R3 Repr. Cat. 1; R61 Repr. Cat. 3; R62 Xn; R20/22 R33 N; R50-53	E; T; N R: 61-3-20/22-33-50/53-62 S: 53-45-60-61		E 1
609-020-00-X	DNOC (ISO); 4,6-dinitro-*o*-cresol	208-601-1	534-52-1	Muta. Cat. 3; R68 T+; R26/27/28 Xi; R38-41 R43 R44 N; R50-53	T+; N R: 26/27/28-38-41-43-44-50/53-68 S: (1/2-)36/37-45-60-61		
609-021-00-5	sodium salt of DNOC; sodium 4,6-dinitro-*o*-cresolate; [1] potassium salt of DNOC; potassium 4,6-dinitro-*o*-cresolate [2]	219-007-7 [1] [2]	2312-76-7 [1] 5787-96-2 [2]	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)13-45-60-61		
609-022-00-0	ammonium salt of DNOC; ammonium 4,6-dinitro-*o*-tolyl oxide	221-037-0	2980-64-5	T+; R26/27/28 R33 N; R50-53	T+; N R: 26/27/28-33-50/53 S: (1/2-)13-28-45-60-61		

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609-023-00-6	dinocap (ISO)	254-408-0	39300-45-3	Repr. Cat. 2; R61 Xn; R20-48/22 Xi; R38 R43 N; R50-53	T; N R: 61-20-22-38-43-48/22-50/53 S: 53-45-60-61		E
609-024-00-1	binapacryl (ISO); 2-*sec*-butyl-4,6-dinitrophenyl-3-methylcrotonate	207-612-9	485-31-4	Repr. Cat. 2; R61 Xn; R21/22 N; R50-53	T; N R: 61-21/22-50/53 S: 53-45-60-61		E
609-025-00-7	dinoseb (ISO); 6-*sec*-butyl-2,4-dinitrophenol	201-861-7	88-85-7	R44 T; R24/25 Repr. Cat. 2; R61 Repr. Cat. 3; R62 Xi; R36 N; R50-53	T; N R: 61-62-24/25-36-44-50/53 S: 53-45-60-61		E
609-026-00-2	salts and esters of dinoseb, with the exception of those specified elsewhere in this Annex	—	—	R44 Repr. Cat. 2; R61 Repr. Cat. 3; R62 T; R24/25 Xi; R36 N; R50-53	T; N R: 61-62-24/25-36-44-50/53 S: 53-45-60-61		AE
609-027-00-8	dinocton; reaction mass of isomers: methyl 2-octyl-4,6-dinitrophenyl carbonate, methyl 4-octyl-2,6-dinitrophenyl carbonate	—	63919-26-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
609-028-00-3	dinex (ISO); 2-cyclohexyl-4,6-dinitrophenol	205-042-5	131-89-5	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		
609-029-00-9	salts and esters of dinex	—	—	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		A

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609-030-00-4	dinoterb (ISO); 2-*tert*-butyl-4,6-dinitrophenol	215-813-8	1420-07-1	Repr. Cat. 2; R61 T+; R28 T; R24 R44 N; R50-53	T+; N R: 61-24-28-44-50/53 S: 53-45-60-61		E
609-031-00-X	salts and esters of dinoterb	—	—	Repr. Cat. 2; R61 T+; R28 T; R24 N; R50-53	T+; N R: 61-24-28-50/53 S: 45-53-60-61		AE
609-032-00-5	bromofenoxim (ISO); 3,5-dibromo-4-hydroxybenzaldehyde-O-(2,4-dinitrophenyl)-oxime	236-129-6	13181-17-4	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)25-60-61		
609-033-00-0	dinosam (ISO); 2-(1-methylbutyl)-4,6-dinitrophenol	—	4097-36-3	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		
609-034-00-6	salts and esters of dinosam	—	—	T; R23/24/25 N; R50-53	T; N R: 23/24/25-50/53 S: (1/2-)13-45-60-61		A
609-035-00-1	nitroethane	201-188-9	79-24-3	R10 Xn; R20/22	Xn R: 10-20/22 S: (2-)9-25-41	Xn; R20/22: C ≥ 12,5 %	
609-036-00-7	nitromethane	200-876-6	75-52-5	R5-10 Xn; R22	Xn R: 5-10-22 S: (2-)41	Xn; R22: C ≥ 12,5 %	
609-037-00-2	5-nitroacenaphthene	210-025-0	602-87-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		
609-038-00-8	2-nitronaphthalene	209-474-5	581-89-5	Carc. Cat. 2; R45 N; R51-53	T; N R: 45-51/53 S: 53-45-61		
609-039-00-3	4-nitrobiphenyl	202-204-7	92-93-3	Carc. Cat. 2; R45 N; R51-53	T; N R: 45-51/53 S: 53-45-61		

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609-040-00-9	nitrofen (ISO); 2,4-dichlorophenyl 4-nitrophenyl ether	217-406-0	1836-75-5	Carc. Cat. 2; R45 Repr. Cat. 2; R61 Xn; R22 N; R50-53	T; N R: 45-61-22-50/53 S: 53-45-60-61		E
609-041-00-4	2,4-dinitrophenol	200-087-7	51-28-5	T; R23/24/25 R33 N; R50	T; N R: 23/24/25-33-50 S: (1/2-)28-37-45-61		
609-042-00-X	pendimethalin (ISO); *N*-(1-ethylpropyl)-2,6-dinitro-3,4-xylidine	254-938-2	40487-42-1	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-29-37-60-61		
609-043-00-5	quintozene (ISO); pentachloronitrobenzene	201-435-0	82-68-8	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)13-24-37-60-61		
609-044-00-0	tecnazene (ISO); 1,2,4,5-tetrachloro-3-nitrobenzene	204-178-2	117-18-0	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
609-045-00-6	reaction mass of: 4,6-dinitro-2-(3-octyl)phenyl methyl carbonate and 4,6-dinitro-2-(4-octyl)phenyl methyl carbonate; dinocton-6	—	8069-76-9	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
609-046-00-1	trifluralin (ISO) (containing < 0.5 ppm NPDA); α,α,α-trifluoro-2,6-dinitro-*N*,*N*-dipropyl-*p*-toluidine (containing < 0.5 ppm NPDA); 2,6-dinitro-*N*,*N*-dipropyl-4-trifluoromethylaniline (containing < 0.5 ppm NPDA); *N*,*N*-dipropyl-2,6-dinitro-4-trifluoromethylaniline (containing < 0.5 ppm NPDA)	216-428-8	1582-09-8	Xi; R36 R43 N; R50-53	Xi; N R: 36-43-50/53 S: (2-)24-37-60-61		
609-047-00-7	2-nitroanisole	202-052-1	91-23-6	Carc. Cat. 2; R45 Xn; R22	T R: 45-22 S: 53-45		E
609-048-00-2	sodium 3-nitrobenzenesulphonate	204-857-3	127-68-4	Xi; R36 R43	Xi R: 36-43 S: (2-)24-26-37		

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609-049-00-8	2,6-dinitrotoluene	210-106-0	606-20-2	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 R52-53	T R: 45-23/24/25-48/22-62-68-52/53 S: 53-45-61		E
609-050-00-3	2,3-dinitrotoluene	210-013-5	602-01-7	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 N; R50-53	T; N R: 45-23/24/25-48/22-62-68-50/53 S: 53-45-60-61		E
609-051-00-9	3,4-dinitrotoluene	210-222-1	610-39-9	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 N; R51-53	T; N R: 45-23/24/25-48/22-62-68-51/53 S: 53-45-61		E
609-052-00-4	3,5-dinitrotoluene	210-566-2	618-85-9	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 R52-53	T R: 45-23/24/25-48/22-62-68-52/53 S: 53-45-61		E
609-053-00-X	hydrazine-trinitromethane	414-850-9	—	E; R3 O; R8 Carc. Cat. 2; R45 T; R23/25 R43	E; T R: 45-3-8-23/25-43 S: 53-45		E

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609-054-00-5	2,3-dinitrophenol; [1] 2,5-dinitrophenol; [2] 2,6-dinitrophenol; [3] 3,4-dinitrophenol; [4] salts of dinitrophenol [5]	200-628-7 [1] 206-348-1 [2] 209-357-9 [3] 209-415-3 [4] [5]	66-56-8 [1] 329-71-5 [2] 573-56-8 [3] 577-71-9 [4] [5]	T; R23/24/25 R33 N; R51-53	T; N R: 23/24/25-33-51 S: (1/2-)28-37-45-61		
609-055-00-0	2,5-dinitrotoluene	210-581-4	619-15-8	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Repr. Cat. 3; R62 T; R23/24/25 Xn; R48/22 N; R51-53	T; N R: 45-23/24/25-48/22-62-68-51/53 S: 53-45-61		E
609-056-00-6	2,2-dibromo-2-nitroethanol	412-380-9	69094-18-4	E; R2 Carc. Cat. 3; R40 Xn; R22-48/22 C; R35 R43 N; R50-53	E; C; N R: 2-22-35-40-43-48/22-50/53 S: (1/2-)23-26-36/37/39-45-60-61	Xn; R22: C ≥ 10 % C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
609-057-00-1	3-chloro-2,4-difluoronitrobenzene	411-980-8	3847-58-3	Xn; R22 C; R34 R43 N; R50-53	C; N R: 22-34-43-50/53 S: (1/2-)22-26-28-36/37/39-45-60-61		
609-058-00-7	2-nitro-2-phenyl-1,3-propanediol	410-360-4	5428-02-4	T; R39-48/25 Xn; R21/22 Xi; R41 R43 N; R51-53	T; N R: 21/22-39-41-43-48/25-51/53 S: 53-45-61		
609-059-00-2	2-chloro-6-(ethylamino)-4-nitrophenol	411-440-1	131657-78-8	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)22-24-37/39-61		

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609-060-00-8	4-[(3-hydroxypropyl)amino]-3-nitrophenol	406-305-9	92952-81-3	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
609-061-00-3	(*E*,*Z*)-4-chlorophenyl(cyclopropyl)ketone *O*-(4-nitrophenylmethyl)oxime	406-100-4	94097-88-8	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
609-062-00-9	2-bromo-2-nitropropanol	407-030-7	24403-04-1	T; R24 Xn; R22-48/22 C; R34 R43 N; R50-53	T; N R: 22-24-34-43-48/22-50/53 S: (1/2-)26-36/37/39-45-60-61		
609-063-00-4	2-[(4-chloro-2-nitrophenyl)amino]ethanol	413-280-8	59320-13-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-61		
609-064-00-X	mesotrione (ISO); 2-[4-(methylsulfonyl)-2-nitrobenzoyl]-1,3-cyclohexanedione	—	104206-82-8	N; R50-53	N R: 50/53 S: 60-61		
609-065-00-5	2-nitrotoluene	201-853-3	88-72-2	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 3; R62 Xn; R22 N; R51-53	T; N R: 45-46-22-62-51/53 S: 53-45-61		E
609-066-00-0	lithium sodium 3-amino-10-{4-(10-amino-6,13-dichloro-4,11-disulfonatobenzo[5,6][1,4]oxazino[2,3-b]phenoxazine-3-ylamino)-6-[methyl(2-sulfonato-ethyl)amino]-1,3,5-triazin-2-ylamino}-6,13-dichlorobenzo[5,6][1,4]oxazino[2,3-b]phenoxazine-4,11-disulfonate	418-870-9	154212-58-5	Xn; R20/21/22-68/20/21/22	Xn R: 20/21/22-68/20/21/22 S: (2-)36/37		
609-067-00-6	sodium and potassium 4-(3-aminopropylamino)-2,6-bis[3-(4-methoxy-2-sulfophenylazo)-4-hydroxy-2-sulfo-7-naphthylamino]-1,3,5-triazine	416-280-6	156769-97-0	R43	Xi R: 43 S: (2-)22-24-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
609-068-00-1	musk xylene; 5-*tert*-butyl-2,4,6-trinitro-*m*-xylene	201-329-4	81-15-2	E; R2 Carc. Cat. 3; R40 N; R50-53	E; Xn; N R: 2-40-50/53 S: (2-)36/37-46-60-61		
609-070-00-2	1,4-dichloro-2-(1,1,2,3,3,3-hexafluoropropoxy)-5-nitrobenzene	415-580-4	130841-23-5	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)36/37/39-60-61		
609-071-00-8	reaction mass of: 2-methylsulfanyl-4,6-bis-(2-hydroxy-4-methoxy-phenyl)-1,3,5-triazine; 2-(4,6-bis-methylsulfanyl-1,3,5-triazin-2-yl)-5-methoxy-phenol	423-520-3	156137-33-6	R43	Xi R: 43 S: (2-)22-24-37		
610-001-00-3	trichloronitromethane; chloropicrin	200-930-9	76-06-2	Xn; R22 T+; R26 Xi; R36/37/38	T+ R: 22-26-36/37/38 S: (1/2-)36/37-38-45		
610-002-00-9	1,1-dichloro-1-nitroethane	209-854-0	594-72-9	T; R23/24/25	T R: 23/24/25 S: (1/2-)26-45		
610-003-00-4	chlorodinitrobenzene	—	—	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)28-36/37-45-60-61		C
610-004-00-X	2-chloro-1,3,5-trinitrobenzene	201-864-3	88-88-0	E; R2 T+; R26/27/28 N; R50-53	E; T+; N R: 2-26/27/28-50/53 S: (1/2-)28-36/37-45-60-61		
610-005-00-5	1-chloro-4-nitrobenzene	202-809-6	100-00-5	Carc. Cat. 3; R40 Muta. Cat. 3; R68 T; R23/24/25 Xn; R48/20/21/22 N; R51-53	T; N R: 23/24/25-40-48/20/21/22-68-51/53 S: (1/2-)28-36/37-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
610-006-00-0	chloronitroanilines with the exception of those specified elsewhere in this Annex	—	—	T+; R26/27/28 R33 N; R51-53	T+; N R: 26/27/28-33-51/53 S: (1/2-)28-36/37-45-61		A C
610-007-00-6	1-chloro-1-nitropropane	209-990-0	600-25-9	Xn; R20/22	Xn R: 20/22 S: (2-)	Xn; R20/22: C ≥ 5 %	
610-008-00-1	2,6-dichloro-4-nitroanisole	403-350-6	17742-69-7	T; R25 N; R51-53	T; N R: 25-51/53 S: (1/2-)36/37-45-61		
610-009-00-7	2-chloro-4-nitroaniline	204-502-2	121-87-9	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-24-61		
610-010-00-2	2-bromo-1-(2-furyl)-2-nitroethylene	406-110-9	35950-52-8	Xn; R22-48/22 C; R34 R43 N; R50-53	C; N R: 22-34-43-48/22-50/53 S: (1/2-)22-26-36/37/39-45-60-61		
611-001-00-6	azobenzene	203-102-5	103-33-3	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R20/22-48/22 N; R50-53	T; N R: 45-20/22-48/22-68-50/53 S: 53-45-60-61		E
611-002-00-1	azoxybenzene	207-802-1	495-48-7	Xn; R20/22	Xn R: 20/22 S: (2-)28		
611-003-00-7	fenaminosulf (ISO); sodium 4-dimethylaminobenzenediazosulphonate	205-419-4	140-56-7	T; R25 Xn; R21 R52-53	T R: 21-25-52/53 S: (1/2-)36/37-45-61		
611-004-00-2	methyl-ONN-azoxymethyl acetate; methyl azoxy methyl acetate	209-765-7	592-62-1	Carc. Cat. 2; R45 Repr. Cat. 2; R61	T R: 45-61 S: 53-45		

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611-005-00-8	disodium {5-[(4'-((2,6-hydroxy-3-((2-hydroxy-5-sulphophenyl)azo)phenyl)azo)(1,1'-biphenyl)-4-yl)azo]salicylato(4-)}cuprate(2-); CI Direct Brown 95	240-221-1	16071-86-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		
611-006-00-3	4-*o*-tolylazo-*o*-toluidine; 4-amino-2',3-dimethylazobenzene; fast garnet GBC base; AAT; *o*-aminoazotoluene	202-591-2	97-56-3	Carc. Cat. 2; R45 R43	T R: 45-43 S: 53-45		
611-007-00-9	tricyclazole (ISO); 5-methyl-1,2,4-triazolo(3,4-b)benzo-1,3-thiazole	255-559-5	41814-78-2	Xn; R22	Xn R: 22 S: (2-)		
611-008-00-4	4-aminoazobenzene; 4-phenylazoaniline	200-453-6	60-09-3	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
611-009-00-X	sodium (1-(5-(4-(4-anilino-3-sulphophenylazo)-2-methyl-5-methylsulphonamidophenylazo)-4-hydroxy-2-oxido-3-(phenylazo)phenylazo)-5-nitro-4-sulphonato-2-naphtholato) iron(II)	401-220-3	—	Xn; R20 R52-53	Xn R: 20-52/53 S: (2-)61		
611-010-00-5	2'-(2-cyano-4,6-dinitrophenylazo)-5'-(*N*,*N*-dipropylamino) propionanilide	403-010-7	106359-94-8	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
611-011-00-0	*N*,*N*,*N*',*N*'-tetramethyl-3,3'-(propylenebis(iminocarbonyl-4,1-phenylenazo(1,6-dihydro-2-hydroxy-4-methyl-6-oxopyridine-3,1-diyl)))di(propylammonium) dilactate	403-340-1	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
611-012-00-6	reaction mass of 2,2-iminodiethanol 6-methyl-2-(4-(2,4,6-triaminopyrimidin-5-ylazo)phenyl)benzothiazole-7-sulfonate and 2-methylaminoethanol 6-methyl-2-(4-(2,4,6-triamino-pyrimidin-5-ylazo)phenyl)benzothiazole-7-sulfonate and *N*,*N*-diethylpropane-1,3-diamine 6-methyl-2-(4-(2,4,6-triamino-pyrimidin-5-ylazo)phenyl)benzothiazole-7-sulfonate	403-410-1	114565-65-0	R43	Xi R: 43 S: (2-)22-24-26-37		
611-013-00-1	trilithium-1-hydroxy-7-(3-sulfonatoanilino)-2-(3-methyl-4-(2-methoxy-4-(3-sulfonatophenylazo)phenylazo)phenylazo) naphthalene-3-sulfonate	403-650-7	117409-78-6	E; R2 N; R51-53	E; N R: 2-51/53 S: (2-)35-61		
611-014-00-7	(tetrasodium 1-(4-(3-acetamido-4-(4'-nitro-2,2'-disulfonatostilben-4-ylazo)anilino)-6-(2,5-disulfonatoanilino)-1,3,5-triazin-2-yl)-3-carboxypyridinium) hydroxide	404-250-5	115099-55-3	R43	Xi R: 43 S: (2-)22-24-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-015-00-2	tetrasodium 4-amino-5-hydroxy-6-(4-(2-(2-(sulfonatooxy) ethylsulfonyl)ethylcarbamoyl)phenylazo)-3-(4-(2-(sulfonatooxy)ethylsulfonyl)phenylazo)naphthalene-2,7-disulfonate	404-320-5	116889-78-2	R43	Xi R: 43 S: (2-)22-24-37		
611-016-00-8	reaction mass of 1,1'-((dihydroxyphenylene)bis(azo-3,1-phenylenazo(1-(3-dimethylaminopropyl)-1,2-dihydro-6-hydroxy-4-methyl-2-oxopyridine-5,3-diyl)))dipyridinium dichloride dihydrochloride, mixed isomers and 1-(1-(3-dimethylaminopropyl)-5-(3-((4-(1-(3-dimethylaminopropyl)-1,6-dihydro-2-hydroxy-4-methyl-6-oxo-5-pyridinio-3-pyridylazo)phenylazo)-2,4(or2,6 or3,5)-dihydroxyphenylazo)phenylazo)-1,2-dihydro-6-hydroxy-4-methyl-2-oxo-3-pyridyl)pyridinium dichloride	404-540-1	—	R43	Xi R: 43 S: (2-)22-24-37		
611-017-00-3	2-(4-(diethylaminopropylcarbamoyl)phenylazo)-3-oxo-*N*-(2,3-dihydro-2-oxobenzimidazol-5-yl)butyramide	404-910-2	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
611-018-00-9	tetraammonium 5-(4-(7-amino-1-hydroxy-3-sulfonato-2-naphthylazo)-6-sulfonato-1-naphthylazo)isophthalate	405-130-5	—	R43	Xi R: 43 S: (2-)24-37		
611-019-00-4	tetralithium 6-amino-4-hydroxy-3-(7-sulfonato-4-(4-sulfonatophenylazo)-1-naphthylazo)naphthalene-2,7-disulfonate	405-150-4	106028-58-4	R43	Xi R: 43 S: (2-)24-37		
611-020-00-X	tetrakis(tetramethylammonium) 6-amino-4-hydroxy-3-(7-sulfonato-4-(4-sulfonatophenylazo)-1-naphthylazo)naphthalene-2,7-disulfonate	405-170-3	116340-05-7	T; R25 R43 R52-53	T R: 25-43-52/53 S: (1/2-)22-24-37-45-61		
611-021-00-5	2-(4-(4-cyano-3-methylisothiazol-5-ylazo)-*N*-ethyl-3-methylanilino)ethyl acetate	405-480-9	—	Xn; R22-48/22 Xi; R38 R53	Xn R: 22-38-48/22-53 S: (2-)22-36/37-61		
611-022-00-0	4-dimethylaminobenzenediazonium 3-carboxy-4-hydroxybenzenesulfonate	404-980-4	—	E; R2 T; R23/25 Xn; R21-48/22 Xi; R41 R43 N; R50/53	E; T; N R: 2-21-23/25-41-43-48/22-50/53 S: (1/2-)3-12-26-35-36/37/39-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-023-00-6	disodium 7-(4,6-dichloro-1,3,5-triazin-2-ylamino)-4-hydroxy-3-(4-(2-(sulfonatooxy)ethylsulfonyl)phenylazo) naphthalene-2-sulfonate	404-600-7	—	R43	X R: 43 S: (2-)22-24-37		
611-024-00-1	Benzidine based azo dyes; 4,4'-diarylazobiphenyl dyes, with the exception of those specified elsewhere in this Annex	—	—	Carc. Cat. 2; R45	T R: 45 S: 53-45		A
611-025-00-7	disodium 4-amino-3-[[4'-[(2,4-diaminophenyl)azo][1,1'-biphenyl]-4-yl]azo]-5-hydroxy-6-(phenylazo)naphtalene-2,7-disulphonate; C.I. Direct Black 38	217-710-3	1937-37-7	Carc. Cat. 2; R45 Repr. Cat. 3; R63	T R: 45-63 S: 53-45		
611-026-00-2	tetrasodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis(azo)]bis[5-amino-4-hydroxynaphthalene-2,7-disulphonate]; C.I. Direct Blue 6	220-012-1	2602-46-2	Carc. Cat. 2; R45 Repr. Cat. 3; R63	T R: 45-63 S: 53-45		
611-027-00-8	disodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis(azo)]bis(4-amino-naphthalene-1-sulphonate); C.I. Direct Red 28	209-358-4	573-58-0	Carc. Cat. 2; R45 Repr. Cat. 3; R63	T R: 45-63 S: 53-45		
611-028-00-3	*C,C'*-azodi(formamide)	204-650-8	123-77-3	R42 R44 ⊗	Xn R: 42-44 S: (2-)22-24-37		
611-029-00-9	*o*-dianisidine based azo dyes; 4,4'-diarylazo-3,3'-dimethoxybiphenyl dyes with the exception of those mentioned elsewhere in this Annex	—	—	Carc. Cat. 2; R45	T R: 45 S: 53-45		A H
611-030-00-4	*o*-tolidine based dyes; 4,4'-diarylazo-3,3'-dimethylbiphenyl dyes, with the exception of those mentioned elsewhere in this Annex	—	—	Carc. Cat. 2; R45	T R: 45 S: 53-45		A H
611-031-00-X	4,4'-(4-iminocyclohexa-2,5-dienylidenemethylene)dianiline hydrochloride; C.I. Basic Red 9	209-321-2	569-61-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		
611-032-00-5	1,4,5,8-tetraaminoanthraquinone; C.I. Disperse Blue 1	219-603-7	2475-45-8	Carc. Cat. 2; R45 Xi; R38-41 R43	T R: 45-38-41-43 S: 53-45		

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611-033-00-0	hexasodium [4,4''-azoxybis(2,2'-disulfonatostilbene-4,4'-diylazo)]-bis[5'-sulfonatobenzene-2,2'- diolato-*O*(2),*O*(2),*N*(1)]-copper(II)	400-020-3	82027-60-9	N; R51-53	N R: 51/53 S: 61		
611-034-00-6	*N*-(5-(bis(2-methoxyethyl)amino)-2-((5-nitro-2,1-benzisothiazol-3-yl)azo)phenylacetamide	402-430-8	105076-77-5	R53	R: 53 S: 61		
611-035-00-1	tetralithium 6-amino-4-hydroxy-3-[7-sulfonato-4-(5-sulfonato-2-naphthylazo)-1-naphthylazo]naphthalene-2,7-disulfonate	403-660-1	107246-80-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
611-036-00-7	2-(4-(5,6(or 6,7)-dichloro-1,3-benzothiazol-2-ylazo)-*N*-methyl-*m*-toluidino)ethyl acetate	405-440-0	—	R43	Xi R: 43 S: (2-)22-24-37		
611-037-00-2	3(or 5)-(4-(*N*-benzyl-*N*-ethylamino)-2-methylphenylazo)-1,4-dimethyl-1,2,4-triazolium methylsulphate	406-055-0	124584-00-5	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)22-24-26-37/39-61		
611-038-00-8	trisodium 1-hydroxynaphthalene-2-azo-4'(5',5''-dimethylbiphenyl)-4''-azo(4''-phenylsulfonyloxybenzene)- 2',2'',4-trisulfonate	406-820-9	—	Xi; R36	Xi R: 36 S: (2-)25-26		
611-039-00-3	7-[((4,6-dichloro-1,3,5-triazin-2-yl)amino)-4-hydroxy-3-(4-((2-sulfoxy)ethyl)sulfonyl)phenylazo]naphthalene-2-sulfonic acid	407-050-6	117715-57-8	R43	Xi R: 43 S: (2-)22-24-37		
611-040-00-9	3-(5-acetylamino-4-(4-[4,6-bis(3-diethylaminopropylamino)-1,3,5-triazin-2-ylamino]phenylazo)-2-(2-methoxyethoxy)phenylazo)-6-amino-4-hydroxy-2-naphthalenesulfonic acid	407-670-7	115099-58-6	N; R50-53	N R: 50/53 S: 60-61		
611-041-00-4	2-[[4[[4,6-bis[[3-(diethylamino)propyl]amino]-1,3,5-triazine-2-yl]amino]phenyl]azo]-*N*-(2,3-dihydro-2-oxo-1*H*-benzimidazol-5-yl)-3-oxobutanamide	407-680-1	98809-11-1	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24-26-37/39-61		
611-042-00-X	trisodium 5-amino-3-[5-(2-bromoacryloylamino)-2-sulfonatophenylazo]-4-hydroxy-6-(4-vinylsulfonylphenylazo) naphthalene-2,7-disulfonate	411-770-6	136213-71-3	R52-53	R: 52/53 S: 61		

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611-043-00-5	reaction mass of: trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6-[2-amino-4-(or 6)-hydroxy-(or 4-amino-2-hydroxy)phenylazo]-6''-(1-carbaniloyl-2-hydroxyprop-1-enylazo)-5',5'''-disulfamoyl-3,3''-disulfonatobis(naphthalene-2,1'-azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate; trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6,6''-bis(1-carbaniloyl-2-hydroxyprop-1-enylazo)-5',5'''-disulfamoyl-3,3''-disulfonatobis(naphthalene-2,1'azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate; trisodium *N*(1')-*N*(2):*N*(1''')-*N*(2'')-η-6,6''-bis[2-amino-4-(or 6)-hydroxy-(or 4-amino-2-hydroxy)phenylazo]5',5'''-disulfamoyl-3,3''-disulfonatobis(naphthalene-2,1'azobenzene-1,2'-diolato-*O*(1),*O*(2'))-chromate (2:1:1)	402-850-1	—	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		

611-044-00-0	reaction mass of: *tert*-alkyl(C_{12}-C_{14})ammonium bis[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]-chromate (1-);
	tert-alkyl(C_{12}-C_{14})ammonium bis[1-[(2-hydroxy-4-nitrophenyl)azo]-2-naphthalenolato(2-)]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium bis[1-[[5-(1,1-dimethylpropyl)-2-hydroxy-3-nitrophenyl]azo]-2-naphthalenolato(2-)]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium [[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]-[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium [[1-[[5-(1,1-dimethylpropyl)-2-hydroxy-3-nitrophenyl]azo]-2-naphthalenolato(2-)]-[1-[(2-hydroxy-5-nitrophenyl)azo]-2-naphthalenolato(2-)]]-chromate(1-);
	tert-alkyl(C_{12}-C_{14})ammonium ((1-(4(or 5)-nitro-2-oxidophenylazo)-2-naphtholato)(1-(3-nitro-2-oxido-5-pentylphenylazo)-2-naphtholato))chromate(1-)
		403-720-7	117527-94-3	N; R51-53	N
					R: 51/53
					S: 61		

611-045-00-6	2-[4-[*N*-(4-acetoxybutyl)-*N*-ethyl]amino-2-methylphenylazo]-3-acetyl-5-nitrothiophene	404-830-8	—	R53	R: 53 S: 61		
611-046-00-1	4,4'-diamino-2-methylazobenzene	407-590-2	43151-99-1	T; R25 Xn; R48/22 R43 N; R50-53	T; N R: 25-43-48/22-50/53 S: (1/2-)22-28-36/37-45-60-61		
611-047-00-7	reaction mass of: 2-[[4-[*N*-ethyl-*N*-(2-acetoxyethyl)amino]phenyl]azo]-5,6-dichlorobenzothiazole; 2-[[4-[*N*-ethyl-*N*-(2-acetoxyethyl)amino]phenyl]azo]-6,7-dichlorobenzothiazole (1:1)	407-890-3	111381-11-4	R53	R: 53 S: 61		

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611-048-00-2	reaction mass of: 2-[[4-[bis(2-acetoxyethyl)amino]phenyl] azo]-5,6-dichlorobenzothiazole; 2-[[4-[bis(2-acetoxyethyl)amino]phenyl]azo]-6,7-dichlorobenzothiazole (1:1)	407-900-6	111381-12-5	R53	R: 53 S: 61		
611-049-00-8	reaction mass of 7-[4-(3-diethylaminopropylamino)-6-(3-diethylammoniopropylamino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(4-phenylazophenylazo)-naphthalene-2-sulfonate, acetic acid, lactic acid (2:1:1)	408-000-6	118658-98-3	Xn; R48/22 R43 R52-53	Xn R: 43-48/22-52/53 S: (2-)22-36/37-61		
611-051-00-9	2-(4-(*N*-ethyl-*N*-(2-hydroxy)ethyl)amino-2-methylphenyl)azo-6-methoxy-3-methyl-benzothiazolium chloride	411-110-7	136213-74-6	N; R50-53	N R: 50/53 S: 60-61		
611-052-00-4	monosodium aqua-[5-[[2,4-dihydroxy-5-[(2-hydroxy-3,5-dinitrophenyl)azo]phenyl]azo]-2-naphthalensulfonate], iron complex	400-720-9	—	R52-53	R: 52/53 S: 61		
611-053-00-X	2,2'-azobis[2-methylpropionamidine] dihydrochloride	221-070-0	2997-92-4	Xn; R22 R43	Xn R: 22-43 S: (2-)24-37		
611-055-00-0	C.I. Disperse Yellow 3; *N*-[4-[(2-hydroxy-5-methylphenyl)azo]phenyl]acetamide	220-600-8	2832-40-8	Carc. Cat. 3; R40 R43	Xn R: 40-43 S: (2-)22-36/37-46		
611-056-00-6	C.I. Solvent Yellow 14; 1-phenylazo-2-naphthol	212-668-2	842-07-9	Carc. Cat. 3; R40 Muta. Cat. 3; R68 R43 R53	Xn R: 40-43-53-68 S: (2-)22-36/37-46-61		
611-057-00-1	6-hydroxy-1-(3-isopropoxypropyl)-4-methyl-2-oxo-5-[4-(phenylazo)phenylazo]-1,2-dihydro-3-pyridinecarbonitrile	400-340-3	85136-74-9	Carc. Cat. 2; R45 R53	T R: 45-53 S: 53-45-61		
611-058-00-7	(6-(4-hydroxy-3-(2-methoxyphenylazo)-2-sulfonato-7-naphthylamino)-1,3,5-triazin-2,4-diyl)bis[(amino-1-methylethyl)ammonium] formate	402-060-7	108225-03-2	Carc. Cat. 2; R45 Xi; R41 N; R51-53	T; N R: 45-41-51/53 S: 53-45-61		
611-059-00-2	octasodium 2-(6-(4-chloro-6-(3-(*N*-methyl-*N*-(4-chloro-6-(3,5-disulfonato-2-naphthylazo)-1-hydroxy-6-naphthylamino)-1,3,5-triazin-2-yl)aminomethyl)phenylamino)-1,3,5-triazin-2-ylamino)-3,5-disulfonato-1-hydroxy-2-naphthylazo) naphthalene-1,5-disulfonate	412-960-1	148878-21-1	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)22-24-26-37/39-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-060-00-8	reaction mass of: sodium 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonatonaphthalen-2-ylazo]-isophthalate; ammonium 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonatonaphthalen-2-ylazo]-isophthalate; 5-[8-[4-[4-[4-[7-(3,5-dicarboxylatophenylazo)-8-hydroxy-3,6-disulfonatonaphthalen-1-ylamino]-6-hydroxy-1,3,5-triazin-2-yl]-2,5-dimethylpiperazin-1-yl]-6-hydroxy-1,3,5-triazin-2-ylamino]-1-hydroxy-3,6-disulfonaphthalen-2-ylazo]-isophthalic acid	413-180-4	187285-15-0	Xi; R41	Xi R: 41 S: (2-)22-26-39		
611-061-00-3	disodium 5-[5-[4-(5-chloro-2,6-difluoropyrimidin-4-ylamino) benzamido]-2-sulfonatophenylazo]-1-ethyl-6-hydroxy-4-methyl-2-oxo-3-pyridylmethylsulfonate	412-530-3	—	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
611-062-00-9	octasodium 2-(8-(4-chloro-6-(3-((4-chloro-6-(3,6-disulfonato-2-(1,5-disulfonatonaphthalen-2-ylazo)-1-hydroxynaphthalen-8-ylamino)-1,3,5-triazin-2-yl)aminomethyl)phenylamino)-1,3,5-triazin-2-ylamino)-3,6-disulfonato-1-hydroxynaphthalen-2-ylazo)naphthalene-1,5-disulfonate	413-550-5	—	Xi; R38-41	Xi R: 38-41 S: (2-)22-26-37/39		
611-063-00-4	trisodium [4'-(8-acetylamino-3,6-disulfonato-2-naphthylazo)-4"-(6-benzoylamino-3-sulfonato-2-naphthylazo)-biphenyl-1,3',3",1'"-tetraolato-*O*,*O*',*O*",*O*"']copper(II)	413-590-3	164058-22-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		
611-064-00-X	4-(3,4-dichlorophenylazo)-2,6-di-*sec*-butyl-phenol	410-600-8	124719-26-2	Xn; R48/22 Xi; R38 N; R50-53	Xn; N R: 38-48/22-50/53 S: (2-)23-25-36/37-60-61		
611-065-00-5	4-(4-nitrophenylazo)-2,6-di-*sec*-butyl-phenol	410-610-2	111850-24-9	Xn; R48/22 Xi; R36/38 R43 N; R50-53	Xn; N R: 36/38-43-48/22-50/53 S: (2-)23-26-36/37-60-61		
611-066-00-0	tetrasodium 5-[4-chloro-6-(*N*-ethyl-anilino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(1,5-disulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate	411-540-5	130201-57-9	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)22-24-26-37/39-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-067-00-6	reaction mass of: bis(tris(2-(2-hydroxy(1-methyl)ethoxy)ethyl) ammonium) 7-anilino-4-hydroxy-3-(2-methoxy-5-methyl-4-(4-sulfonatophenylazo)phenylazo)naphthalene-2-sulfonate; bis(tris(2-(2-hydroxy(2-methyl)ethoxy)ethyl)ammonium) 7-anilino-4-hydroxy-3-(2-methoxy-5-methyl-4-(4-sulfonatophenylazo)phenylazo)naphthalene-2-sulfonate	406-910-8	—	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)26-36/39-61		
611-068-00-1	tetrasodium 4-amino-3,6-bis(5-[4-chloro-6-(2-hydroxyethylamino)-1,3,5-triazin-2-ylamino]-2-sulfonatophenylazo)-5-hydroxynaphthalene-2,7-disulfonate	400-690-7	85665-98-1	N; R51-53	N R: 51/53 S: 61		
611-069-00-7	*N*,*N*-di-[poly(oxyethylene)-co-poly(oxypropylene)]-4-[(3,5-dicyano-4-methyl-2-thienyl)azo)]-3-methylaniline	413-380-1	—	N; R51-53	N R: 51/53 S: 61		
611-070-00-2	reaction mass of: disodium (6-(4-anisidino)-3-sulfonato-2-(3,5-dinitro-2-oxidophenylazo)-1-naphtholato)(1-(5-chloro-2-oxidophenylazo)-2-naphtholato)chromate(1-); trisodium bis(5-(4-anisidino)-3-sulfonato-2-(3,5-dinitro-2-oxidophenylazo)-1-naphtholato)chromate(1-)	405-665-4	—	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
611-071-00-8	tris(tetramethylammonium) 5-hydroxy-1-(4-sulphonatophenyl)-4-(4-sulphonatophenylazo)pyrazole-3-carboxylate	406-073-9	131013-81-5	T; R25 R52-53	T R: 25-52/53 S: (1/2-)37-45-61		
611-072-00-3	2,4-bis[2,2'-[2-(*N*,*N*-dimethylamino)ethyloxycarbonyl]phenylazo]-1,3-dihydroxybenzene, dihydrochloride	407-010-8	118208-02-9	Xn; R22 Xi; R41 N; R51-53	Xn; N R: 22-41-51/53 S: (2-)26-39-61		
611-073-00-9	dimethyl 3,3'-(*N*-(4-(4-bromo-2,6-dicyanophenylazo)-3-hydroxyphenyl)imino)dipropionate	407-310-9	122630-55-1	R53	R: 53 S: 61		
611-074-00-4	reaction mass of: sodium/potassium (3-(4-(5-(5-chloro-2,6-difluoropyrimidin-4-ylamino)-2-methoxy-3-sulfonatophenylazo)-2-oxidophenylazo)-2,5,7-trisulfonato-4-naphtholato) copper(II); sodium/potassium (3-(4-(5-(5-chloro-4,6-difluoropyrimidin-2-ylamino)-2-methoxy-3-sulfonatophenylazo)-2-oxidophenylazo)-2,5,7-trisulfonato-4-naphtholato)copper(II)	407-100-7	—	R43	Xi R: 43 S: (2-)22-24-37		

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611-075-00-X	reaction mass of: tris(3,5,5-trimethylhexylammonium) 4-amino-3-(4-(4-(2-amino-4-hydroxyphenylazo)anilino)-3-sulfonatophenylazo)-5,6-dihydro-5-oxo-6-phenylhydrazono-naphthalene-2,7-disulfonate; tris(3,5,5-trimethylhexylammonium) 4-amino-3-(4-(4-(4-amino-2-hydroxyphenylazo)anilino)-3-sulfonatophenylazo)-5,6-dihydro-5-oxo-6-phenylhydrazononaphthalene-2,7-disulfonate (2:1)	406-000-0	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
611-076-00-5	3-(2,6-dichloro-4-nitrophenylazo)-1-methyl-2-phenylindole	406-280-4	117584-16-4	N; R50-53	N R: 50/53 S: 60-61		
611-077-00-0	dilithium disodium (5,5'-diamino-(μ-4,4'-dihydroxy-1:2-κ-2,O4,O4',-3,3'-[3,3'-dihydroxy-1:2-κ-2-O3,O3'-biphenyl-4,4'-ylenebisazo-1:2-(N3,Ń4-η:Ń3',N4'-η)]-dinaphthalene-2,7-disulfonato(8)))dicuprate(2-)	407-230-4	126637-70-5	Xn; R22 R43	Xn R: 22-43 S: (2-)22-24-37		
611-078-00-6	(2,2'-(3,3'-dioxidobiphenyl-4,4'-diyldiazo)bis(6-(4-(3-(diethylamino)propylamino)-6-(3-(diethylammonio)propylamino)-1,3,5-triazin-2-ylamino)-3-sulfonato-1-naphtholato))dicopper (II) acetate lactate	407-240-9	159604-94-1	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
611-079-00-1	disodium 7-[4-chloro-6-(*N*-ethyl-*o*-toluidino)-1,3,5-triazin-2-ylamino]-4-hydroxy-3-(4-methoxy-2-sulfonatophenylazo)-2-naphthalenesulfonate	410-390-8	147703-64-8	Xi; R41	Xi R: 41 S: (2-)22-26-39		
611-080-00-7	sodium 3-(2-acetamido-4-(4-(2-hydroxybutoxy)phenylazo) phenylazo)benzenesulfonate	410-150-2	147703-65-9	R43	Xi R: 43 S: (2-)22-24-37		
611-081-00-2	tetrasodium [7-(2,5-dihydroxy-KO2-7-sulfonato-6-[4-(2,5,6-trichloro-pyrimidin-4-ylamino)phenylazo]-(*N*1,*N*7-*N*)-1-naphthylazo)-8-hydroxy-KO8-naphthalene-1,3,5-trisulfonato (6-)]cuprate(II)	411-470-5	141048-13-7	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
611-082-00-8	reaction mass of: pentasodium bis(1-(3(or 5)-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato)ferrate(1-); pentasodium [(1-(3-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato)-(5-(4-anilino-3-sulfonatophenylazo)-4-hydroxy-2-oxidophenylazo)-6-nitro-4-sulfonato-2-naphtholato]ferrate(1-)	407-570-3	—	N; R51-53	N R: 51/53 S: 61		
611-083-00-3	reaction mass of: 2-[*N*-ethyl-4-[(5,6-dichlorobenzothiazol-2-yl)azo]-*m*-toludino]ethyl acetate; 2-[*N*-ethyl-4-[(6,7-dichlorobenzothiazol-2-yl)azo]-*m*-toludino]ethyl acetate (1:1)	411-560-4	—	T; R48/25 R43 N; R51-53	T; N R: 43-48/25-51/53 S: (1/2-)22-36/37-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-084-00-9	reaction mass of: *N*-(4-chlorophenyl)-4-(2,5-dichloro-4-(dimethylsulfamoyl)phenylazo)-3-hydroxy-2-naphthalenecarboxamide; N-(4-chlorophenyl)-4-(2,5-dichloro-4-(methylsulfamoyl)phenylazo)-3-hydroxy-2-naphthalenecarboxamide	412-550-2	—	R53	R: 53 S: 61		
611-085-00-4	reaction mass of: 3-cyano-5-(2-cyano-4-nitro-phenylazo)-2-(2-hydroxy-ethylamino)-4-methyl-6-[3-(2-phenoxyethoxy) propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-6-(2-hydroxy-ethylamino)-4-methyl-2-[3-(2-phenoxyethoxy)propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-2-amino-4-methyl-6-[3-(3-hydroxypropoxy)propylamino]pyridine; 3-cyano-5-(2-cyano-4-nitro-phenylazo)-6-amino-4-methyl-2-[3-(3-methoxypropoxy)propylamino]pyridine	411-880-4	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
611-086-00-X	monolithium 5-[[2,4-dihydroxy-5-[(2-hydroxy-3,5-dinitrophenyl)azo]phenyl]azo]-2-naphthalenesulfonate], iron complex, monohydrate	411-360-7	—	R52-53	R: 52/53 S: 61		
611-087-00-5	reaction mass of: 3-((5-cyano-1,6-dihydro-1,4-dimethyl-2-hydroxyl-6-oxo-3-pyridinyl)azo)-benzoyloxy-2-phenoxyethane; 3-((5-cyano-1,6-dihydro-1,4-dimethyl-2-hydroxy-6-oxo-3-pyridinyl)azo)-benzoyloxy-2-ethyloxy-2-(ethylphenol)	411-710-9	—	R53	R: 53 S: 61		
611-088-00-0	reaction mass of: trilithium 4-amino-3-((4-((4-((2-amino-4-hydroxyphenyl)azo)phenyl)amino)-3-sulfophenyl)azo)-5-hydroxy-6-(phenylazo)naphthalene-2,7-disulfonate; trilithium 4-amino-3-((4-((4-((4-amino-2-hydroxyphenyl)azo) phenyl)amino)-3-sulfophenyl)azo)-5-hydroxy-6-(phenylazo) naphthalene-2,7-disulfonate	411-890-9	—	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)22-26-39-61		
611-089-00-6	2-((4-(ethyl-(2-hydroxyethyl)amino)-2-methylphenyl)azo)-6-methoxy-3-methyl-benzothiazolium methylsulfate	411-100-2	136213-73-5	Xn; R48/22 R43 N; R50-53	Xn; N R: 43-48/22-50/53 S: (2-)22-36/37-60-61		
611-090-00-1	2,5-dibutoxy-4-(morpholin-4-yl)benzenediazonium 4-methylbenzenesulfonate	413-290-2	93672-52-7	F; R11 Xn; R22 Xi; R41 R43 R52-53	F; Xn R: 11-22-41-43-52/53 S: (2-)12-22-24-26-37/39-47-61		
611-091-00-7	sodium (1.0-1.95)/lithium (0.05-1) 5-((5-((5-chloro-6-fluoro-pyrimidin-4-yl)amino)-2-sulfonatophenyl)azo)-1,2-dihydro-6-hydroxy-1,4-dimethyl-2-oxo-3-pyridinemethylsulfonate	413-470-0	134595-59-8	R43	Xi R: 43 S: (2-)22-24/25-37		

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611-092-00-2	*tert*-(dodecyl/tetradecyl)-ammonium bis(3-(4-((5-(1,1-dimethyl-propyl)-2-hydroxy-3-nitrophenyl)azo)-3-methyl-5-hydroxy-(1*H*)pyrazol-1-yl)benzenesulfonamidato)chromate	413-210-6	—	N; R51-53	N R: 51/53 S: 61		
611-093-00-8	sodium 2-(4-(4-fluoro-6-(2-sulfo-ethylamino)-[1,3,5]triazin-2-ylamino)-2-ureido-phenylazo)-5-(4-sulfophenylazo)benzene-1-sulfonate	410-770-3	146177-84-6	R43	Xi R: 43 S: (2-)22-24-37		
611-094-00-3	reaction mass of: 2-[2-acetylamino-4-[*N*,*N*-bis[2-ethoxy-carbonyloxy)ethyl]amino]phenylazo]-5,6-dichloro-1,3-benzothiazole; 2-[2-acetylamino-4-[*N*,*N*-bis[2-ethoxy-carbonyloxy)ethyl] amino]phenylazo]-6,7-dichloro-1,3-benzotriazole (1:1)	411-600-0	143145-93-1	R53	R: 53 S: 61		
611-095-00-9	hexasodium 1,1'-[(1-amino-8-hydroxy-3,6-disulfonate-2,7-naphthalenediyl)bis(azo(4-sulfonate-1,3-phenyl)imino[6-[(4-chloro-3-sulfonatophenyl)amino]-1,3,5-triazin-2,4-diyl]]]bis [3-carboxypyridinium] dihydroxide	412-240-7	89797-03-5	N; R51-53	N R: 51/53 S: 22-61		
611-096-00-4	methyl *N*-[3-acetylamino)-4-(2-cyano-4-nitrophenylazo)phenyl]-*N*-[(1-methoxy)acetyl]glycinate	413-040-2	149850-30-6	R43	Xi R: 43 S: (2-)22-24-37		
611-097-00-X	reaction mass of iron complexes of: 1,3-dihydroxy-4-[(5-phenylaminosulfonyl)-2-hydroxyphenylazo]-*n*-(5-amino-sulfonyl-2-hydroxyphenylazo)benzene and: 1,3-dihydroxy-4-[(5-phenylaminosulfonyl)-2-hydroxyphenylazo]-*n*-[4-(4-nitro-2-sulfophenylamino)phenylazo]benzene (*n*=2,5,6)	414-150-3	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
611-098-00-5	tetrakis(tetramethylammonium)3,3'-(6-(2-hydroxyethylamino) 1,3,5-triazine-2,4-diylbisimino(2-methyl-4,1-phenyleneazo)) bisnaphthalene-1,5-disulfonate	405-950-3	131013-83-7	T; R25 R52-53	T R: 25-52/53 S: (1/2-)37-45-61		
611-099-00-0	(methylenebis(4,1-phenylenazo(1-(3-(dimethylamino)propyl)-1,2-dihydro-6-hydroxy-4-methyl-2-oxopyridine-5,3-diyl)))-1,1'-dipyridinium dichloride dihydrochloride	401-500-5	118658-99-4	Carc. Cat. 2; R45 N; R51-53	T; N R: 45-51/53 S: 53-45-61		
611-100-00-4	potassium sodium 3,3'-(3(or4)-methyl-1,2-phenylenebis (imino(6-chloro)-1,3,5-triazine-4,2-diylimino(2-acetamido-5-methoxy)-4,1-phenylenazo)dinaphthalene-1,5-disulfonate	403-810-6	140876-13-7	Xi; R41	Xi R: 41 S: (2-)26-39		
611-101-00-X	2'-(4-chloro-3-cyano-5-formyl-2-thienyl)azo-5'-diethylaminoacetanilide	405-200-5	104366-25-8	R43	Xi R: 43 S: (2-)22-24-37		

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611-103-00-0	trisodium (1-(3-carboxylato-2-oxido-5-sulfonatophenylazo)-5-hydroxy-7-sulfonatonaphthalen-2-amido)nickel(II)	407-110-1	—	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24-26-37/39-61		

611-104-00-6	reaction mass of: trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2(or 4or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2or 6)-(4-(4-nitro-2-sulfonatoanilino)phenylazo)phenolato)ferrate(1-);
	trisodium bis(2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)ferrate(1-);
	trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2(or 4 or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2 or 6)-(4-nitro-2-sulfonatophenylazo)phenolato)ferrate(1-);
	trisodium (2,4(or 2,6 or 4,6)-bis(3,5-dinitro-2-oxidophenylazo)-5-hydroxyphenolato)(2(or 4 or 6)-(3,5-dinitro-2-oxidophenylazo)-5-hydroxy-4(or 2 or 6)-(3-sulfonatophenylazo) phenolato)ferrate(1-);
	disodium 3,3'-(2,4-dihydroxy-1,3(or 1,5 or 3,5)-phenylenediazo)dibenzenesulfonate
		406-870-1	—	R43
				N; R51-53
					Xi; N
					R: 43-51/53
					S: (2-)24-37-61		

611-105-00-1	sodium 4-(4-chloro-6-(*N*-ethylanilino)-1,3,5-triazin-2-ylamino)-2-(1-(2-chlorophenyl)-5-hydroxy-3-methyl-1*H*-pyrazol-4-ylazo)benzenesulfonate	407-800-2	136213-75-7	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
611-106-00-7	hexasodium 4,4'-dihydroxy-3,3'-bis[2-sulfonato-4-(4-sulfonatophenylazo)phenylazo]-7,7'[*p*-phenylenebis[imino(6-chloro-1,3,5-triazine-4,2-diyl)imino]]dinaphthalene-2-sulfonate	410-180-6	157627-99-1	Xi; R41	Xi R: 41 S: (2-)26-39		
611-107-00-2	potassium sodium 4-(4-chloro-6-(3,6-disulfonato-7-(5,8-disulfonato-naphthalen-2-ylazo)-8-hydroxy-naphthalen-1-ylamino)-1,3,5-triazin-2-ylamino)-5-hydroxy-6-(4-(2-sulfatoethanesulfonyl)-phenylazo)-naphthalene-1,7-disulfonate	412-490-7	—	R43	Xi R: 43 S: (2-)22-24-37		
611-108-00-8	disodium 5-((4-((4-chloro-3-sulfonatophenyl)azo)-1-naphthyl) azo)-8-(phenylamino)-1-naphthalenesulfonate	413-600-6	6527-62-4	R52-53	R: 52/53 S: 61		
611-109-00-3	Reaction products of: copper(II) sulfate and tetrasodium 2,4-bis[6-(2-methoxy-5-sulfonatophenylazo)-5-hydroxy-7-sulfonato-2-naphthylamino]-6-(2-hydroxyethylamino)-1,3,5-triazine (2:1)	407-710-3	—	N; R51-53	N R: 51/53 S: 61		
611-110-00-9	tetra-sodium/lithium 4,4'-bis-(8-amino-3,6-disulfonato-1-naphthol-2-ylazo)-3-methylazobenzene	408-210-8	124605-82-9	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-28-37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-111-00-4	disodium 2-[[4-(2-chloroethylsulfonyl)phenyl]-[(2-hydroxy-5-sulfo-3-[3-[2-(2-(sulfooxy)ethylsulfonyl)ethylazo]-4-sulfobenzoato(3-)cuprate(1-)	414-230-8	—	R43	Xi R: 43 S: (2-)22-24-37		
611-112-00-X	tetrasodium 4-hydroxy-5-[4-[3-(2-sulfatoethanesulfonyl)phenylamino]-6-morpholin-4-yl-1,3,5-triazin-2-ylamino]-3-(1-sulfonatonaphthalen-2-ylazo)naphthalene-2,7-disulfonate	413-070-6	—	R43	Xi R: 43 S: (2-)22-24-37		
611-113-00-5	lithium sodium (2-(((5-((2,5-dichlorophenyl)azo)-2-hydroxyphenyl)methylene)amino)benzoato(2-))(2-((4,5-dihydro-3-methyl-5-oxo-1-phenyl-1*H*-pyrazol-4-yl)azo)-5-sulfobenzoato(3-)) chromate(2-)	414-280-0	149626-00-6	N; R51-53	N R: 51/53 S: 24/25-61		
611-114-00-0	lithium sodium (4-((5-chloro-2-hydroxyphenyl)azo)-2,4-dihydro-5-methyl-3*H*-pyrazol-3-onato(2-))(3-((4,5-dihydro-3-methyl-1-(4-methylphenyl)-5-oxo-1*H*-pyrazol-4-yl)azo)-4-hydroxy-5-nitrobenzenesulfonato(3-)) chromate(2-)	414-250-7	149564-66-9	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)22-26-39-61		
611-115-00-6	trilithium bis(4-((4-(diethylamino)-2-hydroxyphenyl)azo)-3-hydroxy-1-naphthalenesulfonato(3-))chromate(3-)	414-290-5	149564-65-8	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
611-116-00-1	reaction mass of: trisodium 5-{4-chloro-6-[2-(2,6-dichloro-5-cyanopyrimidin-4-ylamino)-propylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalene-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(2,6-dichloro-5-cyanopyrimidin-4-ylamino)-1-methyl-ethylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalene-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(4,6-dichloro-5-cyanopyrimidin-2-ylamino)-propylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate; trisodium 5-{4-chloro-6-[2-(4,6-dichloro-5-cyanopyrimidin-2-ylamino)-1-methyl-ethylamino]-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfonatonaphthalen-2-ylazo)-naphthalene-2,7-disulfonate	414-620-8	—	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
611-117-00-7	1,3-bis{6-fluoro-4-[1,5-disulfo-4-(3-aminocarbonyl-1-ethyl-6-hydroxy-4-methyl-pyrid-2-on-5-ylazo)-phenyl-2-ylamino]-1,3,5-triazin-2-ylamino}propane lithium-, sodium salt	415-100-3	149850-29-3	R43	Xi R: 43 S: (2-)22-24-37		
611-118-00-2	sodium 1,2-bis[4-[4-{4-(4-sulfophenylazo)-2-sulfophenylazo}-2-ureido-phenyl-amino]-6-fluoro-1,3,5-triazin-2-ylamino]-propane, sodium salt	413-990-8		R43	Xi R: 43 S: (2-)22-24-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
611-119-00-8	tetrasodium 4-[4-chloro-6-(4-methyl-2-sulfophenylamino)-1,3,5-triazin-2-ylamino]-6-(4,5-dimethyl-2-sulfophenylazo)-5-hydroxynaphthalene-2,7-disulfonate	415-400-4	148878-22-2	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
611-120-00-3	5-{4-[5-amino-2-[4-(2-sulfoxyethylsulfonyl)phenylazo]-4-sulfo-phenylamino]-6-chloro-1,3,5-triazin-2-ylamino}-4-hydroxy-3-(1-sulfo-naphthalen-2-ylazo)-naphthalene-2,7-di-sulfonicacid sodium salt	418-340-7	157707-94-3	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)22-26-39-61		
611-121-00-9	Main component 6 (isomer): asym. 1:2 Cr(III)-complex of: A: 3-hydroxy-4-(2-hydroxy-naphthalene-1-ylazo)naphthalene-1-sulfonic acid, Na-salt and B: 1-[2-hydroxy-5-(4-methoxy-phenylazo)phenylazo]naphthalene-2-ol; Main component 8 (isomer): asym. 1:2 Cr-complex of: A: 3-hydroxy-4-(2-hydroxy-naphthalene-1-ylazo)-naphthalene-1-sulfonic acid, Na-salt and B: 1-[2-hydroxy-5-(4-methoxy-phenylazo)-phenylazo]-naphthalene-2-ol	417-280-9	30785-74-1	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
611-122-00-4	hexasodium (di[*N*-(3-(4-[5-(5-amino-3-methyl-1-phenylpyrazol-4-yl-azo)-2,4-disulfo-anilino]-6-chloro-1,3,5-triazin-2-ylamino)phenyl)-sulfamoyl](di-sulfo)-phthalocyaninato)nickel	417-250-5	151436-99-6	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
611-123-00-X	3-(2,4-bis(4-((5-(4,6-bis(2-aminopropylamino)-1,3,5-triazin-2-ylamino)-4-hydroxy-2,7-disulfonaphthalen-3-yl)azo)phenylamino)-1,3,5-triazin-6-ylamino)propyldiethylammonium lactate	424-310-4	178452-66-9	Xi; R41	Xi R: 41 S: (2-)26-39		
611-124-00-5	reaction mass of: pentasodium 5-amino-3-(5-{4-chloro-6-[4-(2-sulfoxyethoxysulfonato)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo)-6-[5-(2,3-dibromopropionylamino)-2-sulfonatophenylazo]-4-hydroxynaphthalene-2,7-di-sulfonate; pentasodium 5-amino-6-[5-(2-bromoacryloylamino)-2-sulfonatophenylazo]-3-(5-{4-chloro-6-[4-(2-sulfoxyethoxysulfonato)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo)-4-hydroxynaphthalene-2,7-disulfonate; tetrasodium 5-amino-3-[5-{4-chloro-6-[4-(vinylsulfonyl)phenylamino]-1,3,5-triazin-2-ylamino}-2-sulfonatophenylazo]-6-[5-(2,3-dibromopropionylamino)-2-sulfonatophenylazo]-4-hydroxynaphthalene-2,7-disulfonate	424-320-9		Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
611-125-00-0	reaction mass of: Disodium 6-[3-carboxy-4,5-dihydro-5-oxo-4-sulfonatophenyl)pyrazolin-4-yl-azo]-3-[2-oxido-4-(ethen-sulfonyl)-5-methoxyphenylazo]-4-oxidonaphthalene-2-sulfonate copper (II) complex; Disodium 6-[3-carboxy-4,5-dihydro-5-oxo-4-sulfonatophenyl)pyrazolin-4-yl-azo]-3-[2-oxido-4-(2-hydroxyethylsulfonyl)-5-methoxyphenylazo]-4-oxidonaphthalene-2-sulfonate copper (II) complex	423-940-7	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		

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611-126-00-6	2,6-bis-(2-(4-(4-amino-phenylamino)-phenylazo)-1,3-dimethyl-3*H*-imidazolium)-4-dimethylamino-1,3,5-triazine, dichloride	424-120-1	174514-06-8	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
611-127-00-1	pentasodium 4-amino-6-(5-(4-(2-ethyl-phenylamino)-6-(2-sulfatoethanesulfonyl)-1,3,5-triazin-2-ylamino)-2-sulfonato-phenylazo)-5-hydroxy-3-(4-(2-sulfatoethanesulfonyl)phenylazo)naphthalene-2,7-disulfonate	423-790-2	—	R5 Xi; R41 R43 R52-53	Xi R: 5-41-43-52/53 S: (2-)22-26-36/37/39-41-61		
611-128-00-7	*N*,*N*'-bis{6-chloro-4-[6-(4-vinylsulfonylphenylazo)-2,7-disulfonicacid-5-hydroxynapht-4-ylamino]-1,3,5-triazin-2-yl}-*N*-(2-hydroxyethyl)ethane-1,2-diamine, sodium salt	419-500-9	171599-85-2	Xi; R41 R43	Xi R: 41-43 S: (2-)22-24-26-37/39		
611-129-00-2	reaction mass of: 5-[(4-[(7-amino-1-hydroxy-3-sulfo-2-naphthyl)azo]-2,5-diethoxyphenyl)azo]-2-[(3-phosphonophenyl)azo]benzoic acid; 5-[(4-[(7-amino-1-hydroxy-3-sulfo-2-naphthyl)azo]-2,5-diethoxyphenyl)azo]-3-[(3-phosphonophenyl)azo]benzoic acid	418-230-9	163879-69-4	E; R2 Repr. Cat. 3; R62 Xn; R48/22 R43 N; R51-53	E; Xn; N R: 2-43-48/22-62-51/53 S: (2-)26-35-36/37-61		
611-130-00-8	tetra-ammonium 2-[6-[7-(2-carboxylato-phenylazo)-8-hydroxy-3,6-disulfonato-1-naphthylamino]-4-hydroxy-1,3,5-triazin-2-ylamino]benzoate	418-520-5	183130-96-3	Xi; R36 N; R50-53	Xi; N R: 36-50/53 S: (2-)26-39-60-61		
611-131-00-3	2-[2-hydroxy-3-(2-chlorophenyl)carbamoyl-1-naphthylazo]-7-[2-hydroxy-3-(3-methylphenyl)carbamoyl-1-naphthylazo] fluoren-9-one	420-580-2	151798-26-4	Repr. Cat. 2; R61 R53	T R: 61-53 S: 53-45-61		
611-132-00-9	pentasodium bis{7-[4-(1-butyl-5-cyano-1,2-dihydro-2-hydroxy-4-methyl-6-oxo-3-pyridylazo)phenylsulfonylamino]-5'-nitro-3,3'-disulfonatonaphthalene-2-azobenzene-1,2'-diolato} chromate (III)	419-210-2	178452-71-6	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
611-133-00-4	Product by process iron complex of azo dyestuffs obtained by coupling a mixture of diazotized 2-amino-1-hydroxybenzene-4-sulfanilide and 2-amino-1-hydroxybenzene-4-sulfonamide with resorcin, the obtained mixture being subsequently submitted to a second coupling reaction with a mixture of diazotized 3-aminobenzene-1-sulfonic acid (metanilic acid) and 4'-amino-4-nitro-1,1'-diphenylamine-2-sulfonic acid and metallization with ferric chloride, sodium salt	419-260-5	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)26-39-61		
611-134-00-X	trisodium 2-{α[2-hydroxy-3-[4-chloro-6-[4-(2,3-dibromopropionylamino)-2-sulfonatophenylamino]-1,3,5-triazin-2-ylamino]-5-sulfonatophenylazo]-benzylidenehydrazino}-4-sulfonatobenzoate, copper complex	423-770-3	—	Xi; R41 N; R51-53	Xi; N R: 41-51/53 S: (2-)22-26-39-61		

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611-135-00-5	Reaction product of: 2-[[4-amino-2-ureidophenylazo]-5-[(2-(sulfooxy)ethyl)sulfonyl]]benzenesulfonic acid with 2,4,6-trifluoropyrimidine and partial hydrolysis to the corresponding vinylsulfonyl derivative,mixed potassium/sodium salt	424-250-9	—	Xi; R41 R52-53	Xi R: 41-52/53 S: (2-)26-39-61		
611-136-00-0	2-{4-(2-ammoniopropylamino)-6-[4-hydroxy-3-(5-methyl-2-methoxy-4-sulfamoylphenylazo)-2-sulfonatonaphth-7-ylamino]-1,3,5-triazin-2-ylamino}-2-aminopropyl formate	424-260-3	—	Repr. Cat. 3; R62 Xi; R41 N; R51-53	Xn; N R: 41-62-51/53 S: (2-)22-26-36/37/39-61		
611-137-00-6	6-*tert*-butyl-7-chloro-3-tridecyl-7,7a-dihydro-1*H*-pyrazolo[5,1-c]-1,2,4-triazole	419-870-1	159038-16-1	R53	R: 53 S: 61		
611-138-00-1	2-(4-aminophenyl)-6-*tert*-butyl-1*H*-pyrazolo[1,5-b][1,2,4]triazole	415-910-7	152828-25-6	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
611-140-00-2	azafenidin (ISO) 2-(2,4-dichloro-5-prop-2-ynyloxyphenyl)-5,6,7,8-tetrahydro-1,2,4-triazolo[4,3-a]pyridin-3(2*H*)-one	—	68049-83-2	Repr. Cat. 2; R61 Repr. Cat. 3; R62 Xn; R48/22 N; R50-53	T; N R: 61-48/22-62-50/53 S: 53-45-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	E
612-001-00-9	mono-methylamine; [1] di-methylamine; [2] tri-methylamine [3]	200-820-0 [1] 204-697-4 [2] 200-875-0 [3]	74-89-5 [1] 124-40-3 [2] 75-50-3 [3]	F+; R12 Xn; R20 Xi; R37/38-41	F+; Xn R: 12-20-37/38-41 S: (2-)16-26-39	Xn; R20: C ≥ 5 % Xi; R37/38-41: C ≥ 5 % Xi; R36: 0,5 % ≤ C < 5 %	5
612-001-01-6	mono-methylamine ... %; [1] di-methylamine ... %; [2] tri-methylamine ... % [3]	200-820-0 [1] 204-697-4 [2] 200-875-0 [3]	74-89-5 [1] 124-40-3 [2] 75-50-3 [3]	F+; R12 Xn; R20/22 C; R34	F+; C R: 12-20/22-34 S: (1/2-)3-16-26-29-36/37/39-45	Xn; R20/22: C ≥ 15 % C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	B
612-002-00-4	ethylamine	200-834-7	75-04-7	F+; R12 Xi; R36/37	F+; Xi R: 12-36/37 S: (2-)16-26-29		
612-003-00-X	diethylamine	203-716-3	109-89-7	F; R11 Xn; R20/21/22 C; R35	F; C R: 11-20/21/22-35 S: (1/2-)3-16-26-29-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	

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612-004-00-5	triethylamine	204-469-4	121-44-8	F; R11 Xn; R20/21/22 C; R35	F; C R: 11-20/21/22-35 S: (1/2-)3-16-26-29-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
612-005-00-0	butylamine	203-699-2	109-73-9	F; R11 Xn; R20/21/22 C; R35	F; C R: 11-20/21/22-35 S: (1/2-)3-16-26-29-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
612-006-00-6	ethylenediamine; 1,2-diaminoethane	203-468-6	107-15-3	R10 Xn; R21/22 C; R34 R42/43	C R: 10-21/22-34-42/43 S: (1/2-)23-26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/38: 2 % ≤ C < 10 %	
612-007-00-1	2-aminopropane; isopropylamine	200-860-9	75-31-0	F+; R12 Xi; R36/37/38	F+; Xi R: 12-36/37/38 S: (2-)16-26-29		
612-008-00-7	aniline	200-539-3	62-53-3	Carc. Cat. 3; R40 Muta. Cat. 3; R68 T; R23/24/25-48/23/24/25 Xi; R41 R43 N; R50	T; N R: 23/24/25-40-41-43-48/23/24/25-68-50 S: (1/2-)26-27-36/37/39-45-46-61-63	T; R23/24/25: C ≥ 25 % Xn; R20/21/22: 1 % ≤ C < 25 % T; R48/23/24/25: C ≥ 1 % Xn; R48/20/21/22: 0,2 % ≤ C < 1 %	
612-009-00-2	salts of aniline	—	—	Carc. Cat. 3; R40 Muta. Cat. 3; R68 T; R23/24/25-48/23/24/25 Xi; R41 R43 N; R50	T; N R: 23/24/25-40-41-43-48/23/24/25-68-50 S: (1/2-)26-27-36/37/39-45-61-63	T; R23/24/25: C ≥ 25 % Xn; R20/21/22: 1 % ≤ C < 25 % T; R48/23/24/25: C ≥ 1 % Xn; R48/20/21/22: 0,2 % ≤ C < 1 %	A

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612-010-00-8	chloroanilines, with exception of those specified elsewhere in this Annex	—	—	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)28-36/37-45-60-61		C
612-011-00-3	4-nitrosoaniline	211-535-6	659-49-4	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)25-28		
612-012-00-9	*o*-nitroaniline; [1] *m*-nitroaniline; [2] *p*-nitroaniline [3]	201-855-4 [1] 202-729-1 [2] 202-810-1 [3]	88-74-4 [1] 99-09-2 [2] 100-01-6 [3]	T; R23/24/25 R33 R52-53	T R: 23/24/25-33-52/53 S: (1/2-)28-36/37-45-61		C
612-013-00-4	3-aminobenzene sulphonic acid; metanilic acid	204-473-6	121-47-1	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)25-28		
612-014-00-X	sulphanilic acid; 4-aminobenzenesulphonic acid	204-482-5	121-57-3	Xi; R36/38 R43	Xi R: 36/38-43 S: (2-)24-37		
612-015-00-5	*N*-methylaniline	202-870-9	100-61-8	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)28-36/37-45-60-61		
612-016-00-0	*N*,*N*-dimethylaniline	204-493-5	121-69-7	Carc. Cat. 3; R40 T; R23/24/25 N; R51-53	T; N R: 23/24/25-40-51/53 S: (1/2-)28-36/37-45-61		
612-017-00-6	*N*-methyl-*N*,2,4,6-tetranitroaniline; tetryl	207-531-9	479-45-8	E; R2 ⊗ T; R23/24/25 R33	E; T R: 2-23/24/25-33 S: (1/2-)35-45		

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612-018-00-1	bis(2,4,6-trinitrophenyl)amine; hexyl	205-037-8	131-73-7	E; R2 ⊗ T+; R26/27/28 R33 N; R51-53	E; T+; N R: 2-26/27/28-33-51/53 S: (1/2-)35-36-45-61		
612-019-00-7	dipicrylamine, ammonium salt	220-639-0	2844-92-0	E ⊗ R1 T+; R26/27/28 R33 N; R51-53	E; T+; N R: 1-26/27/28-33-51/53 S: (1/2-)28-36/37-45-61		
612-020-00-2	1-naphthylamine	205-138-7	134-32-7	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)24-61		
612-022-00-3	2-naphthylamine	202-080-4	91-59-8	Carc. Cat. 1; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61	Carc. Cat. 1; R45: C ≥ 0,01 %	E
612-023-00-9	phenylhydrazine; [1] phenylhydrazinium chloride; [2] phenylhydrazine hydrochloride; [3] phenylhydrazinium sulphate (2:1) [4]	202-873-5 [1] 200-444-7 [2] 248-259-0 [3] 257-622-2 [4]	100-63-0 [1] 59-88-1 [2] 27140-08-5 [3] 52033-74-6 [4]	Carc. Cat. 2; R45 Muta. Cat. 3; R68 T; R23/24/25-48/23/24/25 Xi; R36/38 R43 N; R50	T; N R: 45-23/24/25-36/38-43-48/23/24/25-68-50 S: 53-45-61		E
612-024-00-4	*m*-toluidine; 3-aminotoluene	203-583-1	108-44-1	T; R23/24/25 R33 N; R50	T; N R: 23/24/25-33-50 S: (1/2-)28-36/37-45-61		
612-025-00-X	nitrotoluidines, with the exception of those specified elsewhere in this Annex	—	—	T; R23/24/25 R33 N; R51-53	T; N R: 23/24/25-33-51/53 S: (1/2-)28-36/37-45-61		C

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-026-00-5	diphenylamine	204-539-4	122-39-4	T; R23/24/25 R33 N; R50-53	T; N R: 23/24/25-33-50/53 S: (1/2-)28-36/37-45-60-61		
612-027-00-0	xylidines with the exception of those specified elsewhere in this Annex; dimethyl anilines with the exception of those specified elsewhere in this Annex	—	—	T; R23/24/25 R33 N; R51-53	T; N R: 23/24/25-33-51/53 S: (1/2-)28-36/37-45-61		C
612-028-00-6	*p*-phenylenediamine	203-404-7	106-50-3	T; R23/24/25 Xi; R36 R43 N; R50-53	T+; N R: 23/24/25-36-43-50/53 S: (1/2-)28-36/37-45-60-61		
612-029-00-1	benzene-1,4-diamine dihydrochloride; *p*-phenylenediamine dihydrochloride	210-834-9	624-18-0	T; R23/24/25 Xi; R36 R43 N; R50-53	T; N R: 23/24/25-36-43-50/53 S: (1/2-)28-36/37-45-60-61		
612-030-00-7	2-methyl-*p*-phenylenediamine sulphate [1]	210-431-8 [1] 228-871-4 [2]	615-50-9 [1] 6369-59-1 [2]	T; R25 Xn; R20/21 R43 N; R51-53	T; N R: 20/21-25-43-51/53 S: (1/2-)24-37-45-61		
612-031-00-2	*N*,*N*-dimethylbenzene-1,3-diamine; [1] 4-amino-*N*,*N*-dimethylaniline; 3-amino-*N*,*N*'-dimethylaniline [2]	220-623-3 [1] 202-807-5 [2]	2836-04-6 [1] 99-98-9 [2]	T; R23/24/25	T R: 23/24/25 S: (1/2-)28-45		C
612-032-00-8	*N*,*N*,*N*',*N*'-tetramethyl-*p*-phenylenediamine	202-831-6	100-22-1	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)28		
612-033-00-3	2-aminophenol	202-431-1	95-55-6	Xn; R20/22 Muta. Cat. 3; R68	Xn R: 20/22-68 S: (2-)28-36/37		
612-034-00-9	2-amino-4,6-dinitrophenol; picramic acid	202-544-6	96-91-3	E; R1 ⊗ Xn; R20/21/22 R52-53	E; Xn R: 1-20/21/22-52/53 S: (2-)35-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-035-00-4	2-methoxyaniline; *o*-anisidine	201-963-1	90-04-0	Carc. Cat. 2; R45 Muta. Cat. 3; R68 T; R23/24/25	T R: 45-23/24/25-68 S: 53-45		E
612-036-00-X	3,3'-dimethoxybenzidine; *o*-dianisidine	204-355-4	119-90-4	Carc. Cat. 2; R45 Xn; R22	T R: 45-22 S: 53-45		E
612-037-00-5	salts of 3,3'-dimethoxybenzidine; salts of *o*-dianisidine	—	—	Carc. Cat. 2; R45 Xn; R22	T R: 45-22 S: 53-45		AE
612-038-00-0	2-nitro-*p*-anisidine; 4-methoxy-2-nitroaniline	202-547-2	96-96-8	T+; R26/27/28 R33 R52-53	T+ R: 26/27/28-33-52/53 S: (1/2-)28-36/37-45-61		
612-039-00-6	2-ethoxyaniline; *o*-phenetidine	202-356-4	94-70-2	T; R23/24/25 R33	T R: 23/24/25-33 S: (1/2-)28-36/37-45		
612-040-00-1	2,4-dinitroaniline	202-553-5	97-02-9	T+; R26/27/28 R33 N; R51-53	T+; N R: 26/27/28-33-51/53 S: (1/2-)28-36/37-45-61		
612-041-00-7	4,4'-bi-*o*-toluidine	204-358-0	119-93-7	Carc. Cat. 2; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61		E
612-042-00-2	benzidine; 1,1'-biphenyl-4,4'-diamine; 4,4'-diaminobiphenyl; biphenyl-4,4'-ylenediamine	202-199-1	92-87-5	Carc. Cat. 1; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61	Carc. Cat. 1; R45: C ≥ 0,01 %	E
612-043-00-8	*N*,*N*'-dimethylbenzidine	—	2810-74-4	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)22-36		

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612-044-00-3	*N,N'*-diacetylbenzidine	210-338-2	613-35-4	Xn; R20/21/22	Xn R: 20/21/22 S: (2-)22-36		
612-046-00-4	allylamine	203-463-9	107-11-9	F; R11 T; R23/24/25 N; R51-53	F; T; N R: 11-23/24/25-51/53 S: (1/2-)9-16-24/25-45-61		
612-047-00-X	benzylamine	202-854-1	100-46-9	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-36/37/39-45		
612-048-00-5	dipropylamine	205-565-9	142-84-7	F; R11 Xn; R20/21/22 C; R35	F; C R: 11-20/21/22-35 S: (1/2-)16-26-36/37/39-45	C; R35: C ≥ 10 % C; R34: 5 % ≤ C < 10 % Xi; R36/37/38: 1 % ≤ C < 5 %	
612-049-00-0	di-*n*-butylamine; [1] di-*sec*-butylamine [2]	203-921-8 [1] 210-937-9 [2]	111-92-2 [1] 626-23-3 [2]	R10 Xn; R20/21/22	Xn R: 10-20/21/22 S: (2-)		
612-050-00-6	cyclohexylamine	203-629-0	108-91-8	R10 Xn; R21/22 C; R34	C R: 10-21/22-34 S: (1/2-)36/37/39-45	C; R34: C ≥ 10 % Xi; R36/38: 2 % ≤ C < 10 %	
612-051-00-1	4,4'-diaminodiphenylmethane; 4,4'-methylenedianiline	202-974-4	101-77-9	Carc. Cat. 2; R45 Muta. Cat. 3; R68 T; R39/23/24/25 Xn; R48/20/21/22 R43 N; R51-53	T; N R: 45-39/23/24/25-43-48/20/21/22-68-51/53 S: 53-45-61		E

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612-052-00-7	(*S*)-*sec*-butylamine; (*S*)-2-aminobutane; [1] (*R*)-*sec*-butylamine; (*R*)-2-aminobutane; [2] *sec*-butylamine; 2-aminobutane [3]	208-164-7 [1] 236-232-6 [2] 237-732-7 [3]	513-49-5 [1] 13250-12-9 [2] 13952-84-6 [3]	F; R11 Xn; R20/22 C; R35 N; R50	F; C; N R: 11-20/22-35-50 S: (1/2-)9-16-26-28-36/37/39-45-61		C
612-053-00-2	*N*-ethylaniline	203-135-5	103-69-5	T; R23/24/25 R33	T R: 23/24/25-33 S: (1/2-)28-37-45		
612-054-00-8	*N*,*N*-diethylaniline	202-088-8	91-66-7	T; R23/24/25 R33 N; R51-53	T; N R: 23/24/25-33-51/53 S: (1/2-)28-37-45-61	T; R23/24/25: C ≥ 5 % Xn; R20/21/22: 1 % ≤ C < 5 %	
612-055-00-3	*N*-methyl-*o*-toluidine; [1] *N*-methyl-*m*-toluidine; [2] *N*-methyl-*p*-toluidine [3]	210-260-9 [1] 211-795-0 [2] 210-769-6 [3]	611-21-2 [1] 696-44-6 [2] 623-08-5 [3]	T; R23/24/25 R33 R52-53	T R: 23/24/25-33-52/53 S: (1/2-)28-36/37-45-61		C
612-056-00-9	*N*,*N*-dimethyl-*p*-toluidine; [1] *N*,*N*-dimethyl-*m*-toluidine; [2] *N*,*N*-dimethyl-*o*-toluidine [3]	202-805-4 [1] 204-495-6 [2] 210-199-8 [3]	99-97-8 [1] 121-72-2 [2] 609-72-3 [3]	T; R23/24/25 R33 R52-53	T R: 23/24/25-33-52/53 S: (1/2-)28-36/37-45-61	T; R23/24/25: C ≥ 5 % Xn; R20/21/22: 1 % ≤ C < 5 %	C
612-057-00-4	piperazine	203-808-3	110-85-0	C; R34 R42/43 R52-53	C R: 34-42/43-52/53 S: (1/2-)22-26-36/37/39-45-61		
612-058-00-X	2,2'-iminodiethylamine; diethylenetriamine	203-865-4	111-40-0	Xn; R21/22 C; R34 R43	C R: 21/22-34-43 S: (1/2-)26-36/37/39-45		
612-059-00-5	3,6-diazaoctanethylenediamin; triethylenetetramine	203-950-6	112-24-3	Xn; R21 C; R34 R43 R52-53	C R: 21-34-43-52/53 S: (1/2-)26-36/37/39-45-61		

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612-060-00-0	3,6,9-triazaundecamethylenediamine; tetraethylenepentamine	203-986-2	112-57-2	Xn; R21/22 C; R34 R43 N; R51-53	C; N R: 21/22-34-43-51/53 S: (1/2-)26-36/37/39-45-61		
612-061-00-6	3-aminopropyldimethylamine; *N*,*N*-dimethyl-1,3-diaminopropane	203-680-9	109-55-7	R10 Xn; R22 C; R34 R43	C R: 10-22-34-43 S: (1/2-)26-36/37/39-45		
612-062-00-1	3-aminopropyldiethylamine; *N*,*N*-diethyl-1,3-diaminopropane	203-236-4	104-78-9	R10 Xn; R21/22 C; R34 R43	C R: 10-21/22-34-43 S: (1/2-)26-36/37/39-45		
612-063-00-7	3,3'-iminodi(propylamine); dipropylenetriamine	200-261-2	56-18-8	T+; R26 T; R24 Xn; R22 C; R35 R43	T+; C R: 22-24-26-35-43 S: (1/2-)26-28-36/37/39-45		
612-064-00-2	3,6,9,12-tetra-azatetradecamethylenediamine; pentacthylenehexamine	223-775-9	4067-16-7	C; R34 R43 N; R50-53	C; N R: 34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-065-00-8	polyethlyenepolyamines with the exception of those specified elsewhere in this Annex	—	—	Xn; R21/22 C; R34 R43 N; R50-53	C; N R: 21/22-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-066-00-3	dicyclohexylamine	202-980-7	101-83-7	Xn; R22 C; R34 N; R50-53	C; N R: 22-34-50/53 S: (1/2-)26-36/37/39-45-60-61	C; R34: C ≥ 10 % Xi; R36/38: 2 % ≤ C < 10 %	
612-067-00-9	3-aminomethyl-3,5,5-trimethylcyclohexylamine	220-666-8	2855-13-2	Xn; R21/22 C; R34 R43 R52-53	C R: 21/22-34-43-52/53 S: (1/2-)26-36/37/39-45-61		

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612-068-00-4	3,3'-dichlorobenzidine; 3,3'-dichlorobiphenyl-4,4'-ylenediamine	202-109-0	91-94-1	Carc. Cat. 2; R45 Xn; R21 R43 N; R50-53	T; N R: 45-21-43-50/53 S: 53-45-60-61		E
612-069-00-X	salts of 3,3'-dichlorobenzidine; salts of 3,3'-dichlorobiphenyl-4,4'-ylenediamine	— — —	— — —	Carc. Cat. 2; R45 Xn; R21 R43 N; R50-53	T; N R: 45-21-43-50/53 S: 53-45-60-61		AE
612-070-00-5	salts of benzidine [	208-519-6 208-520-1 244-236-4 252-984-8	531-85-1 531-86-2 21136-70-9 36341-27-2	Carc. Cat. 1; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		AE
612-071-00-0	salts of 2-naphthylamine	209-030-0 210-313-6	553-00-4 612-52-2	Carc. Cat. 1; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61		AE
612-072-00-6	biphenyl-4-ylamine; xenylamine; 4-aminobiphenyl	202-177-1	92-67-1	Carc. Cat. 1; R45 Xn; R22	T R: 45-22 S: 53-45		E
612-073-00-1	salts of biphenyl-4-ylamine; salts of xenylamine; salts of 4-aminobiphenyl	—	—	Carc. Cat. 1; R45 Xn; R22	T R: 45-22 S: 53-45		AE
612-074-00-7	benzyldimethylamine	203-149-1	103-83-3	R10 Xn; R20/21/22 C; R34 R52-53	C R: 10-20/21/22-34-52/53 S: (1/2-)26-36-45-61		
612-075-00-2	2-aminoethyldimethylamine; 2-dimethylaminoethylamine	203-541-2	108-00-9	F; R11 Xn; R21/22 C; R35	F; C R: 11-21/22-35 S: (1/2-)16-23-26-28-36-45		

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612-076-00-8	ethyldimethylamine	209-940-8	598-56-1	F+; R12 ⊗ Xn; R20/22 C; R34	F+; C R: 12-20/22-34 S: (1/2-)3-16-26-36-45		
612-077-00-3	dimethylnitrosoamine; *N*-nitrosodimethylamine	200-549-8	62-75-9	Carc. Cat. 2; R45 T+; R26 T; R25-48/25 N; R51-53	T+; N R: 45-25-26-48/25-51/53 S: 53-45-61	Carc. Cat. 2; R45: C ≥ 0,001 %	E
612-078-00-9	2,2'-dichloro-4,4'-methylenedianiline; 4,4'-methylene bis(2-chloroaniline)	202-918-9	101-14-4	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		E
612-079-00-4	salts of 2,2'-dichloro-4,4'-methylenedianiline; salts of 4,4'-methylenebis(2-chloroaniline)	—	—	Carc. Cat. 2; R45 Xn; R22 N; R50-53	T; N R: 45-22-50/53 S: 53-45-60-61		AE
612-080-00-X	4-amino-*N*,*N*-diethylaniline; *N*,*N*-diethyl-p-phenylendiamine	202-214-1	93-05-0	T; R25 C; R34	T R: 25-34 S: (1/2-)26-36-45		
612-081-00-5	salts of 4,4'-bi-*o*-toluidine; salts of 3,3'-dimethylbenzidine; salts of *o*-tolidine	210-322-5 265-294-7 277-985-0	612-82-8 64969-36-4 74753-18-7	Carc. Cat. 2; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61		AE
612-082-00-0	thiourea; thiocarbamide	200-543-5	62-56-6	Carc. Cat. 3; R40 Repr. Cat. 3; R63 Xn; R22 N; R51-53	Xn; N R: 22-40-51/53-63 S: (2-)36/37-61		
612-083-00-6	1-methyl-3-nitro-1-nitrosoguanidine	200-730-1	70-25-7	Carc. Cat. 2; R45 Xn; R20 Xi; R36/38 N; R51-53	T; N R: 45-20-36/38-51/53 S: 53-45-61	Carc. Cat. 2; R45: C ≥ 0,01 %	E

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612-084-00-1	dapsone; 4,4'-diamino diphenyl sulfone	201-248-4	80-08-0	Xn; R22	Xn R: 22 S: (2-)22		
612-085-00-7	4,4'-methylenedi-*o*-toluidine	212-658-8	838-88-0	Carc. Cat. 2; R45 Xn; R22 R43 N; R50-53	T; N R: 45-22-43-50/53 S: 53-45-60-61		E
612-086-00-2	amitraz (ISO); *N*,*N*-bis(2,4-xylyliminomethyl) methylamine	251-375-4	33089-61-1	Xn; R22-48/22 R43 N; R50-53	Xn; N R: 22-43-48/22-50/53 S: (2-)22-24-60-36/37-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
612-087-00-8	guazatine (ISO)		108173-90-6	T+; R26 Xn; R21/22 Xi; R37/38-41 N; R50-53	T+; N R: 21/22-26-37/38-41-50/53 S: (1/2-)26-28-36/37/39-38-45-46-60-61-63		
612-088-00-3	simazine (ISO); 6-chloro-*N*,*N'*-diethyl-1,3,5-triazine-2,4-diamine	204-535-2	122-34-9	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-46-60-61		
612-089-00-9	1,5-naphthylenediamine	218-817-8	2243-62-1	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		
612-090-00-4	2,2'-(nitrosoimino)bisethanol	214-237-4	1116-54-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		
612-091-00-X	*o*-toluidine; 2-aminotoluene	202-429-0	95-53-4	Carc. Cat. 2; R45 T; R23/25 Xi; R36 N; R50	T; N R: 45-23/25-36-50 S: 53-45-61		E
612-092-00-5	*N*,*N'*-(2,2-dimethylpropylidene)hexamethylenediamine	401-660-6	1000-78-8	Xi; R38 R43	Xi R: 38-43 S: (2-)24-37		

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612-093-00-0	3,5-dichloro-4-(1,1,2,2-tetrafluoroethoxy)aniline	401-790-3	104147-32-2	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24/25-26-57-60-61		
612-094-00-6	4-(2-chloro-4-trifluoromethyl)phenoxy-2-fluoroaniline hydrochloride	402-190-4	—	T; R48/25 Xn; R22-48/20 Xi; R41 R43 N; R50-53	T; N R: 22-41-43-48/20-48/25-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-095-00-1	benzyl-2-hydroxydodecyldimethylammonium benzoate	402-610-6	113694-52-3	C; R34 Xn; R22 N; R50-53	C; N R: 22-34-50/53 S: (1/2-)26-28-36/37/39-45-60-61		
612-096-00-7	4,4'-carbonimidoylbis[*N*,*N*-dimethylaniline]	207-762-5	492-80-8	Carc. Cat. 3; R40 Xn; R22 Xi; R36 N; R51-53	Xn; N R: 22-36-40-51/53 S: (2-)36/37-61		
612-097-00-2	salts of 4,4'-carbonimidoylbis[*N*,*N*-dimethylaniline]	—	—	Carc. Cat. 3; R40 Xn; R22 Xi; R36 N; R51-53	Xn; N R: 22-36-40-51/53 S: (2-)36/37-61		A
612-098-00-8	nitrosodipropylamine	210-698-0	621-64-7	Carc. Cat. 2; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61	Carc. Cat. 2; R45: C ≥ 0,001 %	E
612-099-00-3	4-methyl-*m*-phenylenediamine; 2,4-toluenediamine	202-453-1	95-80-7	Carc. Cat. 2; R45 T; R25 Xn; R21 Xi; R36 R43 N; R51-53	T; N R: 45-21-25-36-43-51/53 S: 53-45-61		E

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612-100-00-7	propylenediamine	201-155-9	78-90-0	R10 Xn; R21/22 C; R35	C R: 10-21/22-35 S: (1/2-)26-37/39-45		
612-101-00-2	methenamine; hexamethylenetetramine	202-905-8	100-97-0	F; R11 R42/43	F; Xn R: 11-42/43 S: (2-)16-22-24-37		
612-102-00-8	*N*,*N*-bis(3-aminopropyl)methylamine	203-336-8	105-83-9	T; R23/24 Xn; R22 C; R34	T R: 22-23/24-34 S: (1/2-)26-36/37/39-45		
612-103-00-3	*N*,*N*,*N*',*N*'-tetramethylethylenediamine	203-744-6	110-18-9	F; R11 Xn; R20/22 C; R34	F; C R: 11-20/22-34 S: (1/2-)16-26-36/37/39-45		
612-104-00-9	hexamethylenediamine	204-679-6	124-09-4	Xn; R21/22 Xi; R37 C; R34	C R: 21/22-34-37 S: (1/2-)22-26-36/37/39-45		
612-105-00-4	2-piperazin-1-ylethylamine	205-411-0	140-31-8	Xn; R21/22 C; R34 R43 R52-53	C R: 21/22-34-43-52/53 S: (1/2-)26-36/37/39-45-61		
612-106-00-X	2,6-diethylaniline	209-445-7	579-66-8	Xn; R22	Xn R: 22 S: (2-)23-24		
612-107-00-5	1-phenylethylamine; [1] Dl-α-methylbenzylamine [2]	202-706-6 [1] 210-545-8 [2]	98-84-0 [1] 618-36-0 [2]	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-28-36/37/39-45		
612-108-00-0	3-aminopropyltriethoxysilane	213-048-4	919-30-2	Xn; R22 C; R34	C R: 22-34 S: (1/2-)26-36/37/39-45		

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612-109-00-6	bis(2-dimethylaminoethyl)(methyl)amine	221-201-1	3030-47-5	T; R24 Xn; R22 C; R34	T R: 22-24-34 S: (1/2-)26-36/37/39-45		
612-110-00-1	2,2'-dimethyl-4,4'-methylenebis(cyclohexylamine)	229-962-1	6864-37-5	T; R23/24 Xn; R22 C; R35 N; R51-53	T; C; N R: 22-23/24-35-51/53 S: (1/2-)26-36/37/39-45-61		
612-111-00-7	2-methyl-*m*-phenylenediamine; 2,6-toluenediamine	212-513-9	823-40-5	Muta. Cat. 3; R68 Xn; R21/22 R43 N; R51-53	Xn; N R: 21/22-43-51/53-68 S: (2-)24-36/37-61		
612-112-00-2	*p*-anisidine; 4-methoxyaniline	203-254-2	104-94-9	T+; R26/27/28 R33 N; R50	T+; N R: 26/27/28-33-50 S: (1/2-)28-36/37-45-61		
612-113-00-8	6-methyl-2,4-bis(methylthio)phenylene-1,3-diamine	403-240-8	106264-79-3	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
612-114-00-3	*R*,*R*-2-hydroxy-5-(1-hydroxy-2-(4-phenylbut-2-ylamino)ethyl) benzamide hydrogen 2,3-bis(benzoyloxy)succinate	404-390-7	—	F; R11 R43 R52-53	F; Xi R: 11-43-52/53 S: (2-)24-37-61		
612-115-00-9	dimethyldioctadecylammonium hydrogen sulfate	404-050-8	123312-54-9	Xi; R36 R53	Xi R: 36-53 S: (2-)26-39-61		
612-116-00-4	C_{8-18}alkylbis(2-hydroxyethyl)ammonium bis(2-ethylhexyl) phosphate	404-690-8	68132-19-4	T; R23 C; R34 R43 N; R50-53	T; N R: 23-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-117-00-X	C_{12-14}-*tert*-alkylamine, methylphosphonic acid salt	404-750-3	119415-07-5	Xn; R22 C; R34 N; R51-53	C; N R: 22-34-51/53 S: (1/2-)26-28-36/37/39-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-118-00-5	A reaction mass of: (1,3-dioxo-2*H*-benz(de)isoquinolin-2-ylpropyl)hexadecyldimethylammonium 4-toluenesulfonate; (1,3-dioxo-2*H*-benz(de)isoquinolin-2-ylpropyl)hexadecyldimethylammonium bromide	405-080-4	—	Xi; R41 N; R50-53	X; N R: 41-50/53 S: (2-)22-26-39-60-61		
612-119-00-0	benzyldimethyloctadecylammonium 3-nitrobenzenesulfonate	405-330-2	—	Xi; R38-41 N; R50-53	Xi; N R: 38-41-50/53 S: (2-)26-37/39-60-61		
612-120-00-6	aclonifen (ISO); 2-chloro-6-nitro-3-phenoxyaniline	277-704-1	74070-46-5	N; R50-53	N R: 50/53 S: 60-61		
612-121-00-1	amines, polyethylenepoly-; HEPA	268-626-9	68131-73-7	Xn; R21/22 C; R34 R43 N; R50-53	C; N R: 21/22-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-122-00-7	hydroxylamine	232-259-2	7803-49-8	R5 ⊗ Xn; R22-48/22 Xi; R37/38-41 R43 N; R50	Xn; N R: 5-22-37/38-41-43-48/22-50 S: (2-)22-26-36/37/39-61		
612-123-00-2	hydroxylammonium chloride; hydroxylamine hydrochloride; [1] bis(hydroxylammonium) sulphate; hydroxylamine sulphate (2:1); [2] hydroxylammonium hydrogensulphate; hydroxylamine sulphate (1:1) [3]	226-798-2 [1] 233-118-8 [2] 233-154-4 [3]	5470-11-1 [1] 10039-54-0 [2] 10046-00-1 [3]	⊗ Xn; R22-48/22 Xi; R36/38 R43 N; R50	Xn; N R: 22-36/38-43-48/22-50 S: (2-)22-24-37-61		
612-124-00-8	*N*,*N*,*N*-trimethylanilinium chloride	205-319-0	138-24-9	T; R24/25	T R: 24/25 S: (1/2-)25-39-45-53		
612-125-00-3	2-methyl-*p*-phenylenediamine; 2,5-toluenediamine	202-442-1	95-70-5	T; R25 Xn; R20/21 R43 N; R51-53	T; N R: 20/21-25-43-51/53 S: (1/2-)24-37-45-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-126-00-9	toluene-2,4-diammonium sulphate; 4-methyl-m-phenylenediamine sulfate	265-697-8	65321-67-7	Carc. Cat. 2; R45 T; R25 Xn; R21 Xi; R36 R43 N; R51-53	T; N R: 45-21-25-36-43-51/53 S: 53-45-61		E
612-127-00-4	3-aminophenol	209-711-2	591-27-5	Xn; R20/22 N; R51-53	Xn; N R: 20/22-51/53 S: (2-)28-61		
612-128-00-X	4-aminophenol	204-616-2	123-30-8	Muta. Cat. 3; R68 Xn; R20/22 N; R50-53	Xn; N R: 20/22-50/53-68 S: (2-)28-36/37-60-61		
612-129-00-5	diisopropylamine	203-558-5	108-18-9	F; R11 Xn; R20/22 C; R34	F; C R: 11-20/22-34 S: (1/2-)16-26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
612-130-00-0	2,6-diamino-3,5-diethyltoluene; 4,6-diethyl-2-methyl-1,3-benzenediamine; [1] 2,4-diamino-3,5-diethyltoluene; 2,4-diethyl-6-methyl-1,3-benzenediamine; [2] diethylmethylbenzenediamine [3]	218-255-3 [1] 218-256-9 [2] 270-877-4 [3]	2095-01-4 [1] 2095-02-5 [2] 68479-98-1 [3]	Xn; R21/22-48/22 Xi; R36 N; R50-53	Xn; N R: 21/22-36-48/22-50/53 S: (2-)26-28-36/37/39-60-61		C
612-131-00-6	didecyldimethylammonium chloride	230-525-2	7173-51-5	Xn; R22 C; R34	C R: 22-34 S: (2-)26-36/37/39-45		
612-132-00-1	*N,N'*-diphenyl-*p*-phenylenediamine; *N,N'*-diphenyl-1,4-benzenediamine	200-806-4	74-31-7	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
612-133-00-7	(4-ammonio-*m*-tolyl)ethyl(2-hydroxyethyl)ammonium sulphate; 4-(*N*-ethyl-*N*-2-hydroxyethyl)-2-methylphenylenediamine sulphate	247-162-0	25646-77-9	T; R25 Xn; R48/22 R43 N; R50-53	T; N R: 25-43-48/22-50/53 S: (1/2-)24-37-45-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-134-00-2	*N*-(2-(4-amino-*N*-ethyl-*m*-toluidino)ethyl)methanesulphonamide sesquisulphate; 4-(*N*-ethyl-*N*-2-methanesulphonylaminoethyl)-2-methylphenylenediamine sesquisulphate monohydrate	247-161-5	25646-71-3	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
612-135-00-8	*N*-2-naphthylaniline; *N*-phenyl-2-naphthylamine	205-223-9	135-88-6	Carc. Cat. 3; R40 Xi; R36/38 R43 N; R51-53	Xn; N R: 36/38-40-43-51/53 S: (2-)26-36/37-61		
612-136-00-3	*N*-isopropyl-*N*'-phenyl-*p*-phenylenediamine	202-969-7	101-72-4	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61	R43: C ≥ 0,1 %	
612-137-00-9	4-chloroaniline	203-401-0	106-47-8	Carc. Cat. 2; R45 T; R23/24/25 R43 N; R50-53	T; N R: 45-23/24/25-43-50/53 S: 53-45-60-61		E
612-138-00-4	furalaxyl (ISO); methyl *N*-(2,6-dimethylphenyl)-*N*-(2-furylcarbonyl)-Dl-alaninate	260-875-1	57646-30-7	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)36/37/39-61		
612-139-00-X	mefenacet (ISO); 2-(benzothiazol-2-yloxy)-*N*-methyl-*N*-phenylacetamide	277-328-8	73250-68-7	N; R51-53	N R: 51/53 S: 61		
612-140-00-5	quaternary ammonium compounds, benzyl-C_{8-18}-alkyldimethyl, chlorides	264-151-6	63449-41-2	Xn; R21/22 C; R34 N; R50	C; N R: 21/22-34-50 S: (2-)36/37/39-45-61		
612-141-00-0	4,4'-methylenebis(2-ethylaniline); 4,4'-methylenebis(2-ethylbenzeneamine)	243-420-1	19900-65-3	Carc. Cat. 3; R40 Xn; R22 N; R50-53	Xn; N R: 22-40-50/53 S: (2-)36/37-60-61		
612-142-00-6	biphenyl-2-ylamine	201-990-9	90-41-5	Carc. Cat. 3; R40 Xn; R22 R52-53	Xn R: 22-40-52/53 S: (2-)36/37-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-143-00-1	N^5,N^5-diethyltoluene-2,5-diamine monohydrochloride; 4-diethylamino-2-methylaniline monohydrochloride	218-130-3	2051-79-8	T; R25 Xi; R36 R43 N; R50-53	T; N R: 25-36-43-50/53 S: (1/2-)24-26-37-45-60-61		
612-144-00-7	flumetralin (ISO); *N*-(2-chloro-6-fluorobenzyl)-*N*-ethyl-α,α,α-trifluoro-2,6-dinitro-*p*-toluidine	—	62924-70-3	Xi; R36/38 R43 N; R50-53	Xi; N R: 36/38-43-50/53 S: (2-)36/37-60-61		
612-145-00-2	*o*-phenylenediamine	202-430-6	95-54-5	Carc. Cat. 3; R40 Muta. Cat. 3; R68 T; R25 Xn; R20/21 Xi; R36 R43 N; R50-53	T; N R: 20/21-25-36-40-43-50/53-68 S: (1/2-)28-36/37-45-60-61		
612-146-00-8	*o*-phenylenediamine dihydrochloride	210-418-7	615-28-1	Carc. Cat. 3; R40 Muta. Cat. 3; R68 T; R25 Xn; R20/21 Xi; R36 R43 N; R50-53	T; N R: 20/21-25-36-40-43-50/53-68 S: (1/2-)28-36/37-45-60-61		
612-147-00-3	*m*-phenylenediamine	203-584-7	108-45-2	Muta. Cat. 3; R68 T; R23/24/25 Xi; R36 R43 N; R50-53	T; N R: 23/24/25-36-43-50/53-68 S: (1/2-)28-36/37-45-60-61		
612-148-00-9	*m*-phenylenediamine dihydrochloride	208-790-0	541-69-5	Muta. Cat. 3; R68 T; R23/24/25 Xi; R36 R43 N; R50-53	T; N R: 23/24/25-36-43-50/53-68 S: (1/2-)28-36/37-45-60-61		

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612-149-00-4	1,3-diphenylguanidine	203-002-1	102-06-7	Repr. Cat. 3; R62 Xn; R22 Xi; R36/37/38 N; R51-53	Xn; N R: 22-36/37/38-51/53-62 S: (2-)26-36/37/39-61		
612-150-00-X	spiroxamine (ISO); 8-*tert*-butyl-1,4-dioxaspiro[4.5]decan-2-ylmethyl(ethyl)(propyl)amine	—	118134-30-8	Xn; R20/21/22 Xi; R38 R43 N; R50-53	Xn; N R: 20/21/22-38-43-50/53 S: (2-)36/37/39-46-60-61		
612-151-00-5	diaminotoluene, technical product - reaction mass of [2] and [3]; methyl-phenylenediamine; [1] 4-methyl-*m*-phenylene diamine; [2] 2-methyl-*m*-phenylene diamine [3]	246-910-3 [1] 202-453-1 [2] 212-513-9 [3]	25376-45-8 [1] 95-80-7 [2] 823-40-5 [3]	Carc. Cat. 2; R45 T; R25 Xn; R20/21 Xi; R36 R43 N; R51-53	T; N R: 45-20/21-25-36-43-51/53 S: 53-45-61		E
612-152-00-0	*N*,*N*-diethyl-*N*',*N*'-dimethylpropan-1,3-diyl-diamine	406-610-7	62478-82-4	R10 Xn; R20/22-48/20 C; R35 R52-53	C R: 10-20/22-35-48/20-52/53 S: (1/2-)26-28-36/37/39-45-61		
612-153-00-6	4-[*N*-ethyl-*N*-(2-hydroxyethyl)amino]-1-(2-hydroxyethyl)amino-2-nitrobenzene, monohydrochloride	407-020-2	132885-85-9	Xn; R22 R43 R52-53	Xn R: 22-43-52/53 S: (2-)22-24-37-61		
612-154-00-1	6'-(isobutylethylamino)-3'-methyl-2'-phenylamino-spiro[isobenzo-2-oxofuran-7,9'-[9*H*]-xanthene]	410-890-6	95235-29-3	R53	R: 53 S: 61		
612-155-00-7	2'-anilino-6'-((3-ethoxypropyl)ethylamino)-3'-methylspiro(isobenzo-3-oxofuran)-1-(1*H*)-9'-xanthene	411-730-8	93071-94-4	R53	R: 53 S: 61		
612-156-00-2	reaction mass of: trihexadecylmethylammonium chloride; dihexadecyldimethylammonium chloride	405-620-9	—	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		

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612-157-00-8	(Z)-1-benzo[*b*]thien-2-ylethanone oxime hydrochloride	410-780-8	—	Xn; R22-48/22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-48/22-51/53 S: (2-)22-26-36/37/39-61		
612-158-00-3	reaction mass of: bis(5-dodecyl-2-hydroxybenzald-oximate) copper (II) C_{12}-alkyl group is branched; 4-dodecylsalicylaldoxime	410-820-4	—	R53	R: 53 S: 61		
612-159-00-9	Reaction products of: trimethylhexamethylene diamine (a mixture of 2,2,4-trimethyl-1,6-hexanediamine and 2,4,4-trimethyl-1,6-hexanediamine, EINECS listed), Epoxide 8 (mono[(C_{10}-C_{16}-alkyloxy)methyl]oxirane derivatives) and *p*-toluene-sulfonic acid	410-880-1	—	Xn; R22 C; R34 N; R50-53	C; N R: 22-34-50/53 S: (1/2-)23-26-36/37/39-45-60-61		
612-160-00-4	*p*-toluidine; 4-aminotoluene; [1] toluidinium chloride; [2] toluidine sulphate (1:1) [3]	203-403-1 [1] 208-740-8 [2] 208-741-3 [3]	106-49-0 [1] 540-23-8 [2] 540-25-0 [3]	Carc. Cat. 3; R40 T; R23/24/25 Xi; R36 R43 N; R50	T; N R: 23/24/25-36-40-43-50 S: (1/2-)28-36/37-45-61		
612-161-00-X	2,6-xylidine; 2,6-dimethylaniline	201-758-7	87-62-7	Carc. Cat. 3; R40 Xn; R20/21/22 Xi; R37/38 N; R51-53	Xn; N R: 20/21/22-37/38-40-51/53 S: (2-)23-25-36/37-61		
612-162-00-5	dimethyldioctadecylammonium chloride; DODMAC	203-508-2	107-64-2	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)24-26-39-46-60-61		
612-163-00-0	metalaxyl-M (ISO); mefenoxam; (*R*)-2-[(2,6-dimethylphenyl)-methoxyacetylamino]propionic acid methyl ester	—	70630-17-0	Xn; R22 Xi; R41	Xn R: 22-41 S: (2-)26-39-46		
612-164-00-6	2-butyl-2-ethyl-1,5-diaminopentane	412-700-7	137605-95-9	Xn; R21/22-48/22 C; R34 R43 R52-53	C R: 21/22-34-43-48/22-52/53 S: (1/2-)26-36/37/39-45-61		

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612-165-00-1	*N,N'*-diphenyl-*N,N'*-bis(3-methylphenyl)-(1,1'-diphenyl)-4,4'-diamine	413-810-8	65181-78-4	N; R51-53	N R: 51/53 S: 61		
612-166-00-7	reaction mass of: *cis*-(5-ammonium-1,3,3-trimethyl)-cyclohexanemethylammonium phosphate (1:1); *trans*-(5-ammonium-1,3,3-trimethyl)-cyclohexanemethylammonium phosphate (1:1)	411-830-1	114765-88-7	Xi; R41 R43 R52-53	Xi R: 41-43-52/53 S: (2-)24-26-37/39-61		
612-167-00-2	5-acetyl-3-amino-10,11-dihydro-5*H*-dibenz[*b,f*]azepine-hydrochloride	410-490-1	—	Xn; R22-48/22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-48/22-51/53 S: (2-)22-26-36/37/39-61		
612-168-00-8	3,5-dichloro-2,6-difluoropyrdine-4-amine	220-630-1	2840-00-8	Xn; R21/22 N; R51-53	Xn; N R: 21/22-51/53 S: (2-)36/37-61		
612-170-00-9	4-chlorophenyl cyclopropyl ketone *O*-(4-aminobenzyl)oxime	405-260-2	—	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
612-171-00-4	*N,N,N',N'*-tetraglycidyl-4,4'-diamino-3,3'-diethyldiphenylmethane	410-060-3	130728-76-6	Muta. Cat. 3; R68 R43 N; R51-53	Xn; N R: 43-68-51/53 S: (2-)36/37-61		
612-172-00-X	4,4'-methylenebis(*N,N'*-dimethylcyclohexanamine	412-840-9	13474-64-1	Xn; R22-48/22 C; R35 R52-53	C R: 22-35-48/22-52/53 S: (1/2-)26-36/37/39-45-61		
612-173-00-5	lithium 1-amino-4-(4-*tert*-butylanilino)anthraquinone-2-sulfonate	411-140-0	125328-86-1	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)22-26-36/37/39-61		
612-174-00-0	4,4-dimethoxybutylamine	407-690-6	19060-15-2	Xn; R22 C; R34 R43 R52-53	C R: 22-34-43-52/53 S: (1/2-)26-36/37/39-45-61		

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612-175-00-6	2-(*O*-aminooxy)ethylamine dihydrochloride	412-310-7	37866-45-8	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
612-176-00-1	Polymer of 1,3-dibromopropane and *N*,*N*-diethyl-*N*',*N*'-dimethyl-1,3-propanediamine	410-570-6	143747-73-3	N; R50-53	N R: 50/53 S: 60-61		
612-177-00-7	2-naphthylamino-6-sulfomethylamide	412-120-4	104295-55-8	Xn; R48/22 R43 N; R51-53	Xn; N R: 43-48/22-51/53 S: (2-)22-36/37-61		
612-178-00-2	1,4,7,10-tetraazacyclododecane disulfate	412-080-8	112193-77-8	Xn; R22 Xi; R37-41 R52-53	Xn R: 22-37-41-52/53 S: (2-)26-36/37/39-61		
612-179-00-8	1-(2-propenyl)pyridinium chloride	412-740-5	25965-81-5	Xn; R22 R43	Xn R: 22-43 S: (2-)24-37		
612-180-00-3	3-aminobenzylamine	412-230-2	4403-70-7	Xn; R22 C; R34 N; R51-53	C; N R: 22-34-51/53 S: (1/2-)22-26-36/37/39-45-61		
612-181-00-9	2-phenylthioaniline	413-030-8	1134-94-7	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
612-182-00-4	1-ethyl-1-methylmorpholinium bromide	418-210-1	65756-41-4	Muta. Cat. 3; R68	Xn R: 68 S: (2-)36/37		
612-183-00-X	1-ethyl-1-methylpyrrolidinium bromide	418-200-5	69227-51-6	Muta. Cat. 3; R68	Xn R: 68 S: (2-)36/37		
612-184-00-5	6'-(dibutylamino)-3'-methyl-2'-(phenylamino)spiro[isobenzofuran-1(3*H*),9-(9*H*)-xanthen]-3-one	403-830-5	89331-94-2	R52-53	R: 52/53 S: 61		
612-185-00-0	1-[3-[4-((heptadecafluorononyl)oxy)-benzamido]propyl]-*N*,*N*,*N*-trimethylammonium iodide	407-400-8	59493-72-0	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-186-00-6	bis(*N*-(7-hydroxy-8-methyl-5-phenylphenazin-3-ylidene) dimethylammonium) sulfate	406-770-8	149057-64-7	Xn; R48/22 Xi; R41 R43 N; R50-53	Xn; N R: 41-43-48/22-50/53 S: (2-)22-26-36/37/39-60-61		
612-187-00-1	2,3,4-trifluoroaniline	407-170-9	3862-73-5	Xn; R21/22-48/22 Xi; R38-41 N; R51-53	Xn; N R: 21/22-38-41-48/22-51/53 S: (2-)23-26-36/37/39-61		
612-188-00-7	4,4'-(9*H*-fluoren-9-ylidene)bis(2-chloroaniline)	407-560-9	107934-68-9	N; R51-53	N R: 51/53 S: 61		
612-189-00-2	4-amino-2-(aminomethyl)phenol dihydrochloride	412-510-4	135043-64-0	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)22-24-37-60-61		
612-190-00-8	4,4'-methylenebis(2-isopropyl-6-methylaniline)	415-150-6	16298-38-7	Xn; R48/22 N; R51-53	Xn; N R: 48/22-51/53 S: (2-)36-61		
612-191-00-3	Polymer of allylamine hydrochloride	415-050-2	71550-12-4	Xn; R22 R43	Xn R: 22-43 S: (2-)36/37		
612-192-00-9	2-isopropyl-4-(*N*-methyl)aminomethylthiazole	414-800-6	154212-60-9	Xn; R21/22 Xi; R38-41 N; R51-53	Xn; N R: 21/22-38-41-51/53 S: (2-)26-36/37/39-61		
612-193-00-4	3-methylaminomethylphenylamine	414-570-7	18759-96-1	Xn; R21/22 C; R34 R43 N; R50-53	C; N R: 21/22-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
612-194-00-X	2-hydroxy-3-[(2-hydroxyethyl)-[2-(1-oxotetradecyl)amino] ethyl]amino]-*N*,*N*,*N*-trimethyl-1-propanammonium chloride	414-670-0	141890-30-4	Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-50/53 S: (2-)26-39-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
612-195-00-5	bis[tributyl 4-(methylbenzyl)ammonium] 1,5-naphthalenedisulfonate	415-210-1	160236-81-7	Xn; R20/22 Xi; R41 N; R50-53	Xn; N R: 20/22-41-50/53 S: (2-)26-36/39-60-61		
612-196-00-0	4-chloro-*o*-toluidine; [1] 4-chloro-*o*-toluidine hydrochloride [2]	202-441-6 [1] 221-627-8 [2]	95-69-2 [1] 3165-93-3 [2]	Carc. Cat. 2; R45 Muta. Cat. 3; R68 T; R23/24/25 N; R50-53	T; N R: 45-23/24/25-68-50/53 S: 53-45-60-61		E
612-197-00-6	2,4,5-trimethylaniline; [1] 2,4,5-trimethylaniline hydrochloride [2]	205-282-0 [1] [2]	137-17-7 [1] 21436-97-5 [2]	Carc. Cat. 2; R45 T; R23/24/25 N; R51-53	T; N R: 45-23/24/25-51/53 S: 53-45-61		E
612-198-00-1	4,4'-thiodianiline and its salts	205-370-9	139-65-1	Carc. Cat. 2; R45 Xn; R22 N; R51-53	T; N R: 45-22-51/53 S: 53-45-61		E
612-199-00-7	4,4'-oxydianiline and its salts; *p*-aminophenyl ether	202-977-0	101-80-4	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 3; R62 T; R23/24/25 N; R51-53	T; N R: 45-46-23/24/25-62-51/53 S: 53-45-61		E
612-200-00-0	2,4-diaminoanisole; 4-methoxy-*m*-phenylenediamine; [1] 2,4-diaminoanisole sulphate [2]	210-406-1 [1] 254-323-9 [2]	615-05-4 [1] 39156-41-7 [2]	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R22 N; R51-53	T; N R: 45-22-68-51/53 S: 53-45-61		E
612-201-00-6	*N*,*N*,*N*',*N*'-tetramethyl-4,4'-methylendianiline	202-959-2	101-61-1	Carc. Cat. 2; R45 N; R50-53	T; N R: 45-50/53 S: 53-45-60-61		
612-202-00-1	3,4-dichloroaniline	202-448-4	95-76-1	T; R23/24/25 Xi; R41 R43 N; R50-53	T; N R: 23/24/25-41-43-50/53 S: (1/2-)26-36/37/39-45-60-61		

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612-204-00-2	C.I. Basic Violet 3; 4-[4,4'-bis(dimethylamino) benzhydrylidene]cyclohexa-2,5-dien-1-ylidene]dimethylammonium chloride	208-953-6	548-62-9	Carc. Cat. 3; R40 Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-40-41-50/53 S: (2-)26-36/37/39-46-60-61		
612-205-00-8	C.I. Basic Violet 3 with ≥ 0.1 % of Michler's ketone (EC no. 202-027-5)	208-953-6	548-62-9	Carc. Cat. 2; R45 Xn; R22 Xi; R41 N; R50-53	T; N R: 45-22-41-50/53 S: 53-45-60-61		E
612-206-00-3	famoxadone (ISO); 3-anilino-5-methyl-5-(4-phenoxyphenyl)-1,3-oxazolidine-2,4-dione	—	131807-57-3	Xn; R48/22 N; R50-53	Xn; N R: 48/22-50/53 S: (2-)46-60-61		
612-207-00-9	4-ethoxyaniline; *p*-phenetidine	205-855-5	156-43-4	Muta. Cat. 3; R68 Xn; R20/21/22 Xi; R36 R43	Xn R: 20/21/22-36-43-68 S: (2-)36/37-46		
612-209-00-X	6-methoxy-*m*-toluidine; *p*-cresidine	204-419-1	120-71-8	Carc. Cat. 2; R45 Xn; R22	T R: 45-22 S: 53-45		E
612-210-00-5	5-nitro-*o*-toluidine; [1] 5-nitro-*o*-toluidine hydrochloride [2]	202-765-8 [1] 256-960-8 [2]	99-55-8 [1] 51085-52-0 [2]	Carc. Cat. 3; R40 T; R23/24/25 R52-53	T R: 23/24/25-40-52/53 S: (1/2-)36/37-45-61		
612-211-00-0	N-[(benzotriazole-1-yl)methyl)]-4-carboxybenzenesulfonamide	416-470-9	170292-97-4	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
612-212-00-6	2,6-dichloro-4-trifluoromethylaniline	416-430-0	24279-39-8	Xn; R20/22 Xi; R38 R43 N; R50-53	Xn; N R: 20/22-38-43-50/53 S: (2-)24-37-60-61		

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612-213-00-1	isobutylidene-(2-(2-isopropyl-4,4-dimethyloxazolidine-3-yl)-1,1-dimethylethyl)amine	419-850-2	148348-13-4	C; R34 R52-53	C R: 34-52/53 S: (1/2-)23-26-36/37/39-45-61		
612-214-00-7	4-(2,2-diphenylethenyl)-*N*,*N*-di-phenylbenzenamine	421-390-2	89114-90-9	R53	R: 53 S: 61		
612-215-00-2	3-chloro-2-(isopropylthio)aniline	421-700-6	179104-32-6	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
612-217-00-3	1-methoxy-2-propylamine	422-550-4	37143-54-7	F; R11 C; R34 Xn; R22 R52-53	F; C R: 11-22-34-52/53 S: (1/2-)9-26-36/37/39-45-61		
613-001-00-1	ethyleneimine; aziridine	205-793-9	151-56-4	F; R11 Carc. Cat. 2; R45 Muta. Cat. 2; R46 T+; R26/27/28 C; R34 N; R51-53	F; T+; N R: 45-46-11-26/27/28-34-51/53 S: 53-45-61		D E
613-002-00-7	pyridine	203-809-9	110-86-1	F; R11 Xn; R20/21/22	F; Xn R: 11-20/21/22 S: (2-)26-28	Xn; R20/21/22: C ≥ 5 %	
613-003-00-2	1,2,3,4-tetranitrocarbazole	—	6202-15-9	E ⊗ R1 Xn; R20/21/22	E; Xn R: 1-20/21/22 S: (2-)35		
613-004-00-8	crimidine (ISO); 2-chloro-6-methylpyrimidin-4-yldimethylamine	208-622-6	535-89-7	T+; R28	T+ R: 28 S: (1/2-)36/37-45		
613-007-00-4	desmetryne (ISO); 6-isopropylamino-2-methylamino-4-methylthio-1,3,5-triazine	213-800-1	1014-69-3	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		

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613-008-00-X	dazomet (ISO); tetrahydro-3,5-dimethyl-1,3,5-thiadiazine-2-thione	208-576-7	533-74-4	Xn; R22 Xi; R36 N; R50-53	Xn; N R: 22-36-50/53 S: (2-)15-22-24-60-61		
613-009-00-5	2,4,6-trichloro-1,3,5-triazine; cyanuric chloride	203-614-9	108-77-0	R14 T+; R26 Xn; R22 C; R34 R43	T+; R: 14-22-26-34-43 S: (1/2-)26-28-36/37/39-45-46-63	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
613-010-00-0	ametryn (ISO); 2-ethylamino-4-isopropylamino-6-methylthio-1,3,5-triazine	212-634-7	834-12-8	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)36-60-61		
613-011-00-6	amitrole (ISO); 1,2,4-triazol-3-ylamine	200-521-5	61-82-5	Repr. Cat. 3; R63 Xn; R48/22 N; R51-53	Xn; N R: 48/22-63-51/53 S: (2-)13-36/37-61		
613-012-00-1	bentazone (ISO); 3-isopropyl-2,1,3-benzothiadiazine-4-one-2,2-dioxide	246-585-8	25057-89-0	Xn; R22 Xi; R36 R43 R52-53	Xn R: 22-36-43-52/53 S: (2-)24-37-61		
613-013-00-7	cyanazine (ISO); 2-(4-chloro-6-ethylamino-1,3,5-triazine-2-ylamino)-2-methylpropionitrile	244-544-9	21725-46-2	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)37-60-61		
613-014-00-2	ethoxyquin (ISO); 6-ethoxy-1,2-dihydro-2,2,4-trimethylquinoline	202-075-7	91-53-2	Xn; R22	Xn R: 22 S: (2-)24		
613-015-00-8	fenazaflor (ISO); phenyl 5,6-dichloro-2-trifluoromethylbenzimidazole-1-carboxylate	238-134-9	14255-88-0	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
613-016-00-3	fuberidazole (ISO); 2-(2-furyl)benzimidazole	223-404-0	3878-19-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-60-61		
613-017-00-9	bis (8-hydroxyquinolinium) sulphate	205-137-1	134-31-6	Xn; R22	Xn R: 22 S: (2-)36		

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613-018-00-4	morfamquat (ISO); 1,1'-bis(3,5-dimethylmorpholinocarbonylmethyl)-4,4'-bipyridilium ion	—	7411-47-4	Xn; R22 Xi; R36/37/38 R52-53	Xn R: 22-36/37/38-52/53 S: (2-)22-36-61		
613-019-00-X	thioquinox (ISO); 2-thio-1,3-dithiolo(4,5,b)quinoxaline	202-272-8	93-75-4	Xn; R22	Xn R: 22 S: (2-)24		
613-020-00-5	tridemorph (ISO); 2,6-dimethyl-4-tridecylmorpholine	246-347-3	24602-86-6	Repr. Cat. 2; R61 Xn; R20/22 Xi; R38 N; R50-53	T; N R: 61-20/22-38-50/53 S: 53-45-60-61		E
613-021-00-0	dithianon (ISO); 5,10-dihydro-5,10-dioxonaphtho(2,3-b)(1,4)dithiazine-2,3-dicarbonitrile	222-098-6	3347-22-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24-60-61		
613-022-00-6	pyrethrins including cinerins, with the exception of those specified elsewhere in this Annex	—	—	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61		A
613-023-00-1	2-methyl-4-oxo-3-(penta-2,4-dienyl)cyclopent-2-enyl [1*R*-[1α[*S**(*Z*)],3β]]-chrysanthemate; pyrethrin I	204-455-8	121-21-1	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61		
613-024-00-7	2-methyl-4-oxo-3-(penta-2,4-dienyl)cyclopent-2-enyl[1*R*-[1α[*S**(*Z*)](3β)]]-3-(3-methoxy-2-methyl-3-oxoprop-1-enyl)-2,2-dimethylcyclopropanecarboxylate; pyrethrin II	204-462-6	121-29-9	Xn; R20/21/22 N; R50-53	Xn; N R: 20/21/22-50/53 S: (2-)13-60-61		
613-025-00-2	cinerin I; 3-(but-2-enyl)-2-methyl-4-oxocyclopent-2-enyl 2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate	246-948-0	25402-06-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-026-00-8	cinerin II; 3-(but-2-enyl)-2-methyl-4-oxocyclopent-2-enyl 2,2-dimethyl-3-(3-methoxy-2-methyl-3-oxoprop-1-enyl)cyclopropanecarboxylate	204-454-2	121-20-0	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-027-00-3	piperidine	203-813-0	110-89-4	F; R11 T; R23/24 C; R34	F; T R: 11-23/24-34 S: (1/2-)16-26-27-45	T; R23/24: C ≥ 5 % Xn; R20/21: 1 % ≤ C < 5 % C; R34: C ≥ 5 % Xi: R36/38: 1 % ≤ C < 5 %	

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613-028-00-9	morpholine	203-815-1	110-91-8	R10 Xn; R20/21/22 C; R34	C R: 10-20/21/22-34 S: (1/2-)23-36-45	C; R34: C ≥ 10 % Xi; R36/38: 1 % ≤ C < 10 %	
613-029-00-4	dichloro-1,3,5-triazinetrione; dichloroisocyanuric acid	220-487-5	2782-57-2	O; R8 Xn; R22 R31 Xi; R36/37 N; R50-53	O; Xn; N R: 8-22-31-36/37-50/53 S: (2-)8-26-41-60-61		
613-030-00-X	troclosene potassium; [1] troclosene sodium [2]	218-828-8 [1] 220-767-7 [2]	2244-21-5 [1] 2893-78-9 [2]	O; R8 ⊗ Xn; R22 R31 Xi; R36/37 N; R50-53	O; Xn; N R: 8-22-31-36/37-50/53 S: (2-)8-26-41-60-61	Xn; R22: C ≥ 10 % Xi; R36/37: C ≥ 10 % R31: C ≥ 10 %	
613-030-01-7	troclosene sodium, dihydrate	220-767-7	51580-86-0	Xn; R22 R31 Xi; R36/37 N; R50-53	Xn; N R: 22-31-36/37-50/53 S: (2-)8-26-41-60-61		
613-031-00-5	symclosene; trichloroisocyanuric acid; trichloro-1,3,5-triazinetrion	201-782-8	87-90-1	O; R8 Xn; R22 Xi; R36/37 R31 N; R50-53	O; Xn; N R: 8-22-31-36/37-50/53 S: (2-)8-26-41-60-61		
613-032-00-0	methyl-2,3,5,6-tetrachloro-4-pyridylsulphone; 2,3,5,6-tetrachloro-4-(methylsulphonyl)pyridine	236-035-5	13108-52-6	Xn; R21/22 Xi; R36 R43	Xn R: 21/22-36-43 S: (2-)26-28		
613-033-00-6	2-methylaziridine; propyleneimine	200-878-7	75-55-8	F; R11 Carc. Cat. 2; R45 T+; R26/27/28 Xi; R41 N; R51-53	F; T+; N R: 45-11-26/27/28-41-51/53 S: 53-45-61	Carc. Cat. 2; R45: C ≥ 0,01 %	E
613-034-00-1	1,2-dimethylimidazole	217-101-2	1739-84-0	Xn; R22 Xi; R38-41	Xn R: 22-38-41 S: (2-)24-26		

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613-035-00-7	1-methylimidazole	210-484-7	616-47-7	Xn; R21/22 C; R34	C R: 21/22-34 S: (1/2-)26-36-45		
613-036-00-2	2-methylpyridine; 2-picoline	203-643-7	109-06-8	R10 Xn; R20/21/22 Xi; R36/37	Xn R: 10-20/21/22-36/37 S: (2-)26-36		
613-037-00-8	4-methylpyridine; 4-picoline	203-626-4	108-89-4	R10 T; R24 Xn; R20/22 Xi; R36/37/38	T R: 10-20/22-24-36/37/38 S: (1/2-)26-36-45		
613-038-00-3	6-phenyl-1,3,5-triazine-2,4-diyldiamine; 6-phenyl-1,3,5-triazine-2,4-diamine; benzoguanamine	202-095-6	91-76-9	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
613-039-00-9	ethylene thiourea; imidazolidine-2-thione; 2-imidazoline-2-thiol	202-506-9	96-45-7	Repr. Cat. 2; R61 Xn; R22	T R: 61-22 S: 53-45		E
613-040-00-4	azaconazole (ISO); 1-{[2-(2,4-dichlorophenyl)-1,3-dioxolan-2-yl]methyl}-1*H*-1,2.4-triazole	262-102-3	60207-31-0	Xn; R22	Xn R: 22 S: (2-)46		
613-041-00-X	morpholine-4-carbonyl chloride	239-213-0	15159-40-7	R14 Carc. Cat. 3; R40 Xi; R36/38	Xn R: 14-36/38-40 S: (2-)26-30-36-38		
613-042-00-5	imazalil (ISO); 1-[2-(allyloxy)-2-(2,4-dichlorophenyl)ethyl]-1*H*-imidazole	252-615-0	35554-44-0	Xn; R20/22 Xi; R41 N; R50-53	Xn; N R: 20/22-41-50/53 S: (2-)26-39-60-61		
613-043-00-0	imazalil sulphate (ISO) powder; 1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate; [1] (±)-1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1H-imidazolium hydrogen sulphate [2]	261-351-5 [1] 281-291-3 [2]	58594-72-2 [1] 83918-57-4 [2]	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24/25-37-46-60-61		

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613-043-01-8	imazalil sulphate (ISO), aqueous solution; 1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate; [1] (±)-1- [2-(allyloxy)ethyl-2-(2,4-dichlorophenyl)]-1*H*-imidazolium hydrogen sulphate [2]	261-351-5 [1] 281-291-3 [2]	58594-72-2 [1] 83918-57-4 [2]	Xn; R22 C; R34 R43 N; R50-53	C; N R: 22-34-43-50/53 S: (2-)26-36/37/39-45-60-61	C; R34: C ≥ 50 % Xi; R38: 30 % ≤ C < 50 % Xi; R41: 15 % ≤ C < 50 % Xi; R36: 5 % ≤ C < 15 %	
613-044-00-6	captan (ISO); 1,2,3,6-tetrahydro-*N*-(trichloromethylthio)phthalimide	205-087-0	133-06-2	Carc. Cat. 3; R40 T; R23 Xi; R41 R43 N; R50	T; N R: 23-40-41-43-50 S: (1/2-)26-29-36/37/39-45-61		
613-045-00-1	folpet (ISO); *N*-(trichloromethylthio)phthalimide	205-088-6	133-07-3	Carc. Cat. 3; R40 Xn; R20 Xi; R36 R43 N; R50	Xn; N R: 20-36-40-43-50 S: (2-)36/37-46-61		
613-046-00-7	captafol (ISO); 1,2,3,6-tetrahydro-*N*-(1,1,2,2-tetrachloroethylthio)phthalimide	219-363-3	2425-06-1	Carc. Cat. 2; R45 R43 N; R50-53	T; N R: 45-43-50/53 S: 53-45-60-61		
613-047-00-2	1-dimethylcarbamoyl-5-methylpyrazol-3-yl dimethylcarbamate; dimetilan (ISO)	211-420-0	644-64-4	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)36/37-45-60-61		
613-048-00-8	carbendazim (ISO); methyl benzimidazol-2-ylcarbamate	234-232-0	10605-21-7	Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 N; R50-53	T; N R: 46-60-61-50/53 S: 53-45-60-61		
613-049-00-3	benomyl (ISO); methyl 1-(butylcarbamoyl)benzimidazol-2-ylcarbamate	241-775-7	17804-35-2	Muta. Cat. 2; R46 Repr. Cat. 2; R60-61 Xi; R37/38 R43 N; R50-53	T; N R: 46-60-61-37/38-43-50/53 S: 53-45-60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	

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613-050-00-9	carbadox (INN); methyl 3-(quinoxalin-2-ylmethylene)carbazate 1,4-dioxide; 2-(methoxycarbonylhydrazonomethyl)quinoxaline 1,4-di-oxide	229-879-0	6804-07-5	F; R11 Carc. Cat. 2; R45 Xn; R22	F; T R: 45-11-22 S: 53-45		E
613-051-00-4	molinate (ISO); *S*-ethyl 1-perhydroazepinecarbothioate; *S*-ethyl perhydroazepine-1-carbothioate	218-661-0	2212-67-1	Carc. Cat. 3; R40 Repr. Cat. 3; R62 Xn; R20/2248/22 R43 N; R50-53	Xn; N R: 20/22-40-43-48/22-62-50/53 S: (2-)36/37-46-60-61	N; R50-53: C ≥ 0,25 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
613-052-00-X	trifenmorph (ISO); 4-tritylmorpholine	215-812-2	1420-06-0	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-053-00-5	anilazine (ISO); 2-chloro-*N*-(4,6-dichloro-1,3,5-triazin-2-yl)aniline	202-910-5	101-05-3	Xi; R36/38 N; R50-53	Xi; N R: 36/38-50/53 S: (2-)22-60-61		
613-054-00-0	thiabendazol (ISO); 2-(thiazole-4-yl)benzimidazole	205-725-8	148-79-8	N; R50-53	N R: 50/53 S: 60-61		
613-056-00-1	1,2-dimethyl-3,5-diphenylpyrazolium methylsulphate; difenzoquat methyl sulfate	256-152-5	43222-48-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-057-00-7	dodemorph (ISO); 4-cyclododecyl-2,6-dimethylmorpholine	216-474-9	1593-77-7	Xi; R36/37/38 N; R51-53	Xi; N R: 36/37/38-51/53 S: (2-)26-61		
613-058-00-2	permethrin (ISO); *m*-phenoxybenzyl 3-(2,2-dichlorovinyl)-2,2-dimethylcyclo-propanecarboxylate	258-067-9	52645-53-1	Xn; R20/22 R43 N; R50-53	Xn; N R: 20/22-43-50/53 S: (2-)13-24-36/37/39-60-61	N; R50-53: C ≥ 0,025 % N; R51-53: 0,0025 % ≤ C < 0,025 % R52-53: 0,00025 % ≤ C < 0,0025 %	
613-059-00-8	profluralin (ISO); *N*-(cyclopropylmethyl)-α,α,α-trifluoro-2,6-dinitro-*N*-propyl-*p*-toluidine	247-656-6	26399-36-0	Xi; R36 N; R50-53	Xi; N R: 36-50/53 S: (2-)60-61		

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613-060-00-3	resmethrin (ISO); 5-benzyl-3-furylmethyl (±)-*cis-trans*-chrysanthemate	233-940-7	10453-86-8	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60/61		
613-061-00-9	6-(1α,5aβ,8aβ,9-pentahydroxy-7β-isopropyl-2β,5β,8β-trimethylperhydro-8bα,9-epoxy-5,8-ethanocyclopenta[1,2-b] indenyl) pyrrole-2-carboxylate; ryania	239-732-2	15662-33-6	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
613-062-00-4	sabadilla (ISO); veratrine	—	8051-02-3	Xi; R36/37/38	Xi R: 36/37/38 S: (2-)36/37/39		
613-063-00-X	secbumeton (ISO); 2-sec-butylamino-4-ethylamino-6-methoxy-1,3,5-triazine	247-554-1	26259-45-0	Xn; R22 Xi; R36 N; R50-53	Xn; N R: 22-36-50/53 S: (2-)60-61		
613-064-00-5	5-(3,6,9-trioxa-2-undecyloxy)benzo(d)-1,3-dioxolane; sesamex	—	51-14-9	Xn; R22	Xn R: 22 S: (2-)		
613-065-00-0	simetryn (ISO); 2,4-bis(ethylamino)-6-methylthio-1,3,5-triazine	213-801-7	1014-70-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-066-00-6	terbumeton (ISO); 2-*tert*-butylamino-4-ethylamino-6-methoxy-1,3,5-triazine	251-637-8	33693-04-8	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-067-00-1	propazine (ISO); 2-chloro-4,6-bis(isopropylamino)-1,3,5-triazine	205-359-9	139-40-2	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		
613-068-00-7	atrazine (ISO); 2-chloro-4-ethylamine-6-isopropylamine-1,3,5-triazine	217-617-8	1912-24-9	Xn; R48/22 R43 N; R50-53	Xn; N R: 43-48/22-50/53 S: (2-)36/37-60-61		
613-069-00-2	ε-caprolactam	203-313-2	105-60-2	Xn; R20/22 Xi; R36/37/38	Xn R: 20/22-36/37/38 S: (2-)		

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613-070-00-8	propylenethiourea	—	2122-19-2	Repr. Cat. 3; R63 Xn; R22 R52-53	Xn R: 22-52/53-63 S: (2-)36/37-46-61		
613-071-00-3	2-fluoro-5-trifluoromethylpyridine	400-290-2	69045-82-5	R10 R43 R52-53	Xi R: 10-43-52/53 S: (2-)24-37-61		
613-072-00-9	*N*,*N*-bis(2-ethylhexyl)-((1,2,4-triazol-1-yl)methyl)amine	401-280-0	91273-04-0	C; R34 R43 N; R51-53	C; N R: 34-43-51/53 S: (1/2-)26-28-36/37/39-45-61		
613-073-00-4	*N*,*N*-dimethyl-2-(3-(4-chlorophenyl)-4,5-dihydropyrazol-1-ylphenylsulphonyl)ethylamine	401-410-6	10357-99-0	Xn; R48/22 R43 N; R51-53	Xn; N R: 43-48/22-51/53 S: (2-)24-37-61		
613-074-00-X	3-(3-methylpent-3-yl)isoxazol-5-ylamine	401-460-9	82560-06-3	T; R23/25 Xi; R41 R52-53	T R: 23/25-41-52/53 S: (1/2-)22-26-36/37/39-45-61		
613-075-00-5	1,3-dichloro-5-ethyl-5-methylimidazolidine-2,4-dione	401-570-7	89415-87-2	O; R8 T; R23 C; R34 Xn; R22 R43 N; R50	O; T; N R: 8-22-23-34-43-50 S: (1/2-)8-26-36/37/39-45-61		
613-076-00-0	3-chloro-5-trifluoromethyl-2-pyridylamine	401-670-0	79456-26-1	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
613-077-00-6	reaction mass of 5-heptyl-1,2,4-triazol-3-ylamine and 5-nonyl-1,2,4-triazol-3-ylamine	401-940-8	—	Xn; R22 Xi; R36 N; R51-53	Xn; N R: 22-36-51/53 S: (2-)22-26-61		
613-078-00-1	*N*,*N*,*N*,*N*-tetrakis(4,6-bis(butyl-(*N*-methyl-2,2,6,6-tetramethylpiperidin-4-yl)amino)triazin-2-yl)-4,7-diazadecane-1,10-diamine	401-990-0	106990-43-6	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		

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613-079-00-7	4-(1(or 4 or 5 or 6)-methyl-8,9,10-trinorborn-5-en-2-yl) pyridine, reaction mass of isomers	402-520-7	—	Xn; R21/22 Xi; R38 R43 N; R50-53	Xn; N R: 21/22-38-43-50/53 S: (2-)36/37-60-61		
613-080-00-2	3-(bis(2-ethylhexyl)aminomethyl)benzothiazole-2(3*H*)-thione	402-540-6	105254-85-1	C; R34 R43 N; R50-53	C; N R: 34-43-50/53 S: (1/2-)26-28-36/37/39-45-60-61		
613-081-00-8	1-butyl-2-methylpyridinium bromide	402-680-8	26576-84-1	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
613-082-00-3	2-methyl-1-pentylpyridinium bromide	402-690-2	—	Xn; R21/22 R52-53	Xn R: 21/22-52/53 S: (2-)36/37-61		
613-083-00-9	2-(4-(3-(4-chlorophenyl)-2-pyrazolin-1-yl)phenylsulfonyl) ethyldimethylammonium formate	402-120-2	—	C; R34 Xn; R48/22 R43 N; R50-53	C; N R: 34-43-48/22-50/53 S: (1/2-)24-26-28-37/39-45-60-61		
613-084-00-4	2-(4-(3-(4-chlorophenyl)-4,5-dihydropyrazolyl)phenylsulphonyl)ethyldimethylammonium hydrogen phosphonate	402-490-5	106359-93-7	Xi; R36 N; R50-53	Xi; N R: 36-50/53 S: (2-)26-60-61		
613-085-00-X	reaction mass of 1,1'-(methylenebis(4,1-phenylene))dipyrrole-2,5-dione and *N*-(4-(4-(2,5-dioxopyrrol-1-yl)benzyl)phenyl) acetamide and 1-(4-(4-(5-oxo-2*H*-2-furylidenamino)benzyl) phenyl)pyrrole-2,5-dione	401-970-1	—	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
613-086-00-5	caffeine	200-362-1	58-08-2	Xn; R22	Xn R: 22 S: (2-)		
613-087-00-0	tetrahydrothiophene	203-728-9	110-01-0	F; R11 Xn; R20/21/22 Xi; R36/38 R52-53	F; Xn R: 11-20/21/22-36/38-52/53 S: (2-)16-23-36/37-61		

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613-088-00-6	1,2-benzisothiazol-3(2*H*)-one; 1,2-benzisothiazolin-3-one	220-120-9	2634-33-5	Xn; R22 Xi; R38-41 R43 N; R50	Xn; N R: 22-38-41-43-50 S: (2-)24-26-37/39-61	R43: C ≥ 0,05 %	
613-089-00-1	diquat dibromide; [1] diquat dichloride; [2] 6,7-dihydrodipyrido[1,2-α:2',1'-c]pyrazinediylium dihydroxide [3]	201-579-4 [1] 223-714-6 [2] 301-467-6 [3]	85-00-7 [1] 4032-26-2 [2] 94021-76-8 [3]	T+; R26 T; R48/25 Xn; R22 Xi; R36/37/38 R43 N; R50-53	T+; N R: 22-26-36/37/38-43-48/25-50/53 S: (1/2-)28-36/37/39-45-60-61		
613-090-00-7	paraquat dichloride; 1,1-dimethyl-4,4'-bipyridinium dichloride; [1] paraquat dimethylsulfate; 1,1-dimethyl-4,4'-bipyridinium dimethyl sulphate [2]	217-615-7 [1] 218-196-3 [2]	1910-42-5 [1] 2074-50-2 [2]	T+; R26 T; R24/25-48/25 Xi; R36/37/38 N; R50-53	T+; N R: 24/25-26-36/37/38-48/25-50/53 S: (1/2-)22-28-36/37/39-45-60-61		
613-091-00-2	morfamquat dichloride; [1] morfamquat sulfate [2]	225-062-8 [1] [2]	4636-83-3 [1] 29873-36-7 [2]	Xn; R22 Xi; R36/37/38 R52-53	Xn R: 22-36/37/38-52/53 S: (2-)22-36-61		
613-092-00-8	1,10-phenanthroline	200-629-2	66-71-7	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)45-60-61		
613-093-00-3	hexasodium 6,13-dichloro-3,10-bis((4-(2,5-disulfonatoanilino)-6-fluoro-1,3,5-triazin-2-ylamino)prop-3-ylamino)-5,12-dioxa-7,14-diazapentacene-4,11-disulfonate	400-050-7	85153-92-0	R42/43	Xn R: 42/43 S: (2-)22-24-37		
613-094-00-9	4-methoxy-*N*,6-dimethyl-1,3,5-triazin-2-ylamine	401-360-5	5248-39-5	Xn; R22-48/22	Xn R: 22-48/22 S: (2-)22-36		
613-095-00-4	sodium 3-(2*H*-benzotriazol-2-yl)-5-*sec*-butyl-4-hydroxybenzenesulfonate	403-080-9	92484-48-5	Xi; R41	Xi R: 41 S: (2-)26-39		
613-096-00-X	2-amino-6-ethoxy-4-methylamino-1,3,5-triazine	403-580-7	62096-63-3	Xn; R22	Xn R: 22 S: (2-)		

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613-097-00-5	7-amino-3-((5-carboxymethyl-4-methyl-1,3-thiazol-2-ylthio) methyl)-8-oxo-5-thia-1-azabicyclo(4.2.0)oct-2-ene-2-carboxylic acid	403-690-5	111298-82-9	R42/43 R52-53	Xn R: 42/43-52/53 S: (2-)22-24-37-61		
613-098-00-0	*N*-(*n*-octyl)-2-pyrrolidone	403-700-8	2687-94-7	C; R34 N; R51-53	C; N R: 34-51/53 S: (1/2-)23-26-36/37/39-45		
613-099-00-6	1-dodecyl-2-pyrrolidone	403-730-1	2687-96-9	C; R34 R43 N; R50-53	C; N R: 34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
613-100-00-X	2,9-bis(3-(diethylamino)propylsulfamoyl)quino(2,3-b)acridine-7,14-dione	404-230-6	—	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
613-101-00-5	*N*-*tert*-pentyl-2-benzothiazolesulfenamide	404-380-2	110799-28-5	R43 R52-53	Xi R: 43-52/53 S: (2-)36/37-61		
613-102-00-0	dimethomorph (ISO); 4-(3-(4-chlorophenyl)-3-(3,4-dimethoxyphenyl)acryloyl)morpholine	404-200-2	110488-70-5	N; R51-53	N R: 51/53 S: 61		
613-103-00-6	sodium 5-*n*-butylbenzotriazole	404-450-2	118685-34-0	Xn; R22 C; R34 R43 N; R51-53	C; N R: 22-34-43-51/53 S: (1/2-)26-36/37/39-45-61		
613-104-00-1	5-*tert*-butyl-3-isoxazolylamine hydrochloride	404-840-2	—	Xn; R22-48/22 Xi; R41 R52-53	Xn R: 22-41-48/22-52/53 S: (2-)26-36/39-61		
613-105-00-7	hexakis(tetramethylammonium) 4,4'-vinylenebis((3-sulfonato-4,1-phenylene)imino(6-morpholino-1,3,5-triazine-4,2-diyl) imino)bis(5-hydroxy-6-phenylazonaphthalene-2,7-disulfonate)	405-160-9	124537-30-0	T; R25 R43 R52-53	T R: 25-43-52/53 S: (1/2-)24-37-45-61		
613-106-00-2	tetrapotassium 2-(4-(5-(1-(2,5-disulfonatophenyl)-3-ethoxycarbonyl-5-hydroxypyrazol-4-yl)penta-2,4-dienylidene)-3-ethoxycarbonyl-5-oxo-2-pyrazolin-1-yl)benzene-1,4-disulfonate	405-240-3	—	R43	Xi R: 43 S: (2-)24-37		

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613-107-00-8	hexasodium 2,2'-vinylenebis((3-sulfonato-4,1-phenylene) imino(6-(*N*-cyanoethyl-*N*-(2-hydroxypropyl)amino)-1,3,5-triazine-4,2-diyl)imino)dibenzene-1,4-disulfonate	405-280-1	76508-02-6	Xi; R36	Xi R: 36 S: (2-)26		
613-108-00-3	benzothiazole-2-thiol	205-736-8	149-30-4	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
613-109-00-9	bis(piperidinothiocarbonyl) disulphide	202-328-1	94-37-1	Xi; R36/37/38 R43	Xi R: 36/37/38-43 S: (2-)24-26-37		
613-110-00-4	dimepiperate (ISO); *S*-(1-methyl-1-phenylethyl) piperidine-1-carbothioate	262-784-2	61432-55-1	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
613-111-00-X	1,2,4-triazole	206-022-9	288-88-0	Repr. Cat. 3; R63 Xn; R22 Xi; R36	Xn R: 22-36-63 S: (2-)36/37		
613-112-00-5	octhilinone (ISO); 2-octyl-2*H*-isothiazol-3-one	247-761-7	26530-20-1	T; R23/24 Xn; R22 C; R34 R43 N; R50-53	T; N R: 22-23/24-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61	R43: C ≥ 0,05 %	
613-113-00-0	2-(morpholinothio)benzothiazole	203-052-4	102-77-2	Xi; R36/38 R43 N; R51-53	Xi; N R: 36/38-43-51/53 S: (2-)24-26-37-61		
613-114-00-6	2,2',2"-(hexahydro-1,3,5-triazine-1,3,5-triyl)triethanol; 1,3,5-tris(2-hydroxyethyl)hexahydro-1,3,5-triazine	225-208-0	4719-04-4	Xn; R22 R43	Xn R: 22-43 S: (2-)24-37	R43: C ≥ 0,1 %	
613-115-00-1	hymexazol (ISO); 3-hydroxy-5-methylisoxazole	233-000-6	10004-44-1	Xn; R22 Xi; R41 R52-53	Xn R: 22-41-52/53 S: (2-)26-39-61		

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613-116-00-7	tolylfluanid (ISO); dichloro-*N*-[(dimethylamino)sulphonyl]fluoro-*N*-(*p*-tolyl) methanesulphenamide	211-986-9	731-27-1	T; R23 Xn; R48/20 Xi; R36/37/38 R43 N; R50-53	T; N R: 23-36/37/38-43-48/20-50/53 S: (1/2-)24-26-37-38-45-60-61		
613-117-00-2	diniconazole (ISO); (*E*)-β-[(2,4-dichlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazol-1-ethanol; (*E*)-(*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-en-3-ol	—	76714-88-0	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-118-00-8	flubenzimine (ISO); *N*-[3-phenyl-4,5-bis[(trifluoromethyl)imino]thiazolidin-2-ylidene]aniline;	253-703-1	37893-02-0	Xi; R36 N; R50-53	Xi; N R: 36-50/53 S: (2-)26-60-61		
613-119-00-3	(benzothiazol-2-ylthio)methyl thiocyanate; TCMTB	244-445-0	21564-17-0	T+; R26 Xn; R22 Xi; R36/38 R43 N; R50-53	T+; N R: 22-26-36/38-43-50/53 S: (1/2-)28-36/37-38-45-60-61		
613-120-00-9	bioresmethrin; (5-bezylfur-3-yl)methyl(1R)-*trans*-2,2-dimethyl-3-(2-methylpropenyl)cyclopropanecarboylate	249-014-0	28434-01-7	N; R50-53	N R: 50/53 S: 60-61		
613-121-00-4	chlorsulfuron (ISO); 2-chloro-*N*-[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino] carbonyl]benzenesulphonamide;	265-268-5	64902-72-3	N; R50-53	N R: 50/53 S: 60-61		
613-122-00-X	diclobutrazole (ISO); (*R**, *R**)-(±)-β-[(2,4-dichlorophenyl)methyl]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol; (2*RS*, 3*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1yl)pentan-3-ol	—	75736-33-3	Xi; R36 N; R51-53	Xi; N R: 36-51/53 S: (2-)26-61		
613-123-00-5	5,6-dihydro-3*H*-imidazo[2,1-*c*]-1,2,4-dithiazole-3-thione; etem	251-684-4	33813-20-6	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		

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613-124-00-0	fenpropimorph (ISO); *cis*-4-[3-(*p*-*tert*-butylphenyl)-2-methylpropyl]-2,6-dimethylmorpholine	266-719-9	67564-91-4	Repr. Cat. 3; R63 Xn; R22 Xi; R38 N; R51-53	Xn; N R: 22-38-63-51/53 S: (2-)36/37-46-61		
613-125-00-6	hexythiazox (ISO); *trans*-5-(4-chlorophenyl)-*N*-cyclohexyl-4-methyl-2-oxo-3-thiazolidine-carboxamide	—	78587-05-0	N; R50-53	N R: 50/53 S: 60-61		
613-126-00-1	imazapyr (ISO); 2-[4,5-dihydro-4-methyl-4-(1-methylethyl)-5-oxo-1*H*-imidazol-2-yl]-3-pyridine carboxylate	—	81334-34-1	Xi; R36 R52-53	Xi R: 36-52/53 S: (2-)26-61		
613-127-00-7	1,1-dimethylpiperidinium chloride; mepiquat chloride	246-147-6	24307-26-4	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
613-128-00-2	prochloraz (ISO); *N*-propyl-*N*-[2-(2,4,6-trichlorophenoxy)ethyl]-1*H*-imidazole-1-carboxamide;	266-994-5	67747-09-5	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
613-129-00-8	metamitron (ISO); 4-amino-3-methyl-6-phenyl-1,2,4-triazin-5-one	255-349-3	41394-05-2	Xn; R22 N; R50	Xn; N R: 22-50 S: (2-)61		
613-131-00-9	pyroquilon (ISO); 1,2,5,6-tetrahydropyrrolo[3,2,1-ij]quinolin-4-one	—	57369-32-1	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
613-132-00-4	hexazinone (ISO); 3-cyclohexyl-6-dimethylamino-1-methyl-1,2,3,4-tetrahydro-1,3,5-triazine-2,4-dione;	257-074-4	51235-04-2	Xn; R22 Xi; R36 N; R50-53	Xn; N R: 22-36-50/53 S: (2-)60-61		
613-133-00-X	etridiazole (ISO); 5-ethoxy-3-trichloromethyl-1,2,4-thiadiazole	219-991-8	2593-15-9	Carc. Cat. 3; R40 T; R23 Xn; R21/22 N; R50-53	T; N R: 21/22-23-40-50/53 S: (1/2-)36/37-38-45-60-61		

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613-134-00-5	myclobutanil (ISO); 2-(4-chlorophenyl)-2-(1*H*-1,2,4-triazol-1-ylmethyl)hexanenitrile	—	88671-89-0	Repr. Cat. 3; R63 Xn; R22 Xi; R36 N; R51-53	Xn; N R: 22-36-51/53-63 S: (2-)36/37-46-61		
613-135-00-0	di(benzothiazol-2-yl) disulphide	204-424-9	120-78-5	R31 R43 N; R50-53	Xi; N R: 31-43-50/53 S: (2-)36/37-60-61		
613-136-00-6	*N*-cyclohexylbenzothiazole-2-sulphenamide	202-411-2	95-33-0	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
613-137-00-1	methabenzthiazuron (ISO); 1-(1,3-benzothiazol-2-yl)1,3-dimethylurea	242-505-0	18691-97-9	N; R50-53	N R: 50/53 S: 60-61		
613-138-00-7	quinoxyfen (ISO); 5,7-dichloro-4-(4-fluorophenoxy)quinoline	—	124495-18-7	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-46-60-61		
613-139-00-2	metsulfuron-methyl; 2-(4-methoxy-6-methyl-1,3,5-triazin-2-ylcarbamoylsulfamoyl) benzoic acid	—	74223-64-6	N; R50-53	N R: 50/53 S: 60-61		
613-140-00-8	cycloheximide (ISO); 4-{(2*R*)-2-[(1*S*,3*S*,5*S*)-3,5-dimethyl-2-oxocyclohexyl]-2-hydroxyethyl}piperidine-2,6-dione	200-636-0	66-81-9	Muta. Cat. 3; R68 Repr. Cat. 2; R61 T+; R28 N; R51-53	T+; N R: 61-28-51/53-68 S: 53-45-61		E
613-141-00-3	1,4-diamino-2-(2-butyltetrazol-5-yl)-3-cyanoanthraquinone	401-470-3	93686-63-6	R53	R: 53 S: 61		
613-142-00-9	*trans*-*N*-methyl-2-styryl-[4'-aminomethine-(1-acetyl-1-(2-methoxyphenyl)acetamido)]pyridinium acetate	405-860-4	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)22-24-37-61		
613-143-00-4	1-(3-phenylpropyl)-2-methylpyridinium bromide	405-930-4	10551-42-5	Xn; R22 Xi; R36 R52-53	Xn R: 22-36-52/53 S: (2-)26-36/37-61		

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613-144-00-X	Reaction products of: poly(vinyl acetate), partially hydrolyzed, with (*E*)-2-(4-formylstyryl)-3,4-dimethylthiazoliummethyl sulfate	406-460-2	125139-08-4	R52-53	R: 52/53 S: 61		
613-145-00-5	(*S*)-3-benzyloxycarbonyl-1,2,3,4-tetrahydro-isoquinolinium 4-methylbenzenesulfonate	406-960-0	77497-97-3	N; R51-53	N R: 51/53 S: 61		
613-146-00-0	*N*-ethyl-*N*-methylpiperidinium iodide	407-780-5	4186-71-4	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)22-61		
613-147-00-6	4-[2-(1-methyl-2-(4-morpholinyl)ethoxy)ethyl]morpholine	407-940-4	111681-72-2	Xi; R41	Xi R: 41 S: (2-)26-39		
613-148-00-1	tetrasodium 1,2-bis(4-fluoro-6-[5-(1-amino-2-sulfonatoanthrachinon-4-ylamino)-2,4,6-trimethyl-3-sulfonatophenylamino]-1,3,5-triazin-2-ylamino)ethane	411-240-4	143683-23-2	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24/25-37-61		
613-149-00-7	pyridaben (ISO); 2-*tert*-butyl-5-(4-*tert*-butylbenzylthio)-4-chloropyridazin-3(2*H*)-one	405-700-3	96489-71-3	T; R23/25 N; R50-53	T; N R: 23/25-50/53 S: (1/2-)36/37-45-60-61		
613-150-00-2	2,2'-[3,3'-(piperazine-1,4-diyl)dipropyl]bis(1*H*-benzimidazo[2,1-*b*]benzo[*l*,*m*,*n*][3,8]phenanthroline-1,3,6-trione	406-295-6	—	R53	R: 53 S: 61		
613-151-00-8	1-(3-mesyloxy-5-trityloxymethyl-2-D-threofuryl)thymine	406-360-9	104218-44-2	R53	R: 53 S: 61		
613-152-00-3	phenyl *N*-(4,6-dimethoxypyrimidin-2-yl)carbamate	406-600-2	89392-03-0	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
613-153-00-9	2,3,5-trichloropyridine	407-270-2	16063-70-0	R52-53	R: 52/53 S: 61		
613-154-00-4	2-amino-4-chloro-6-methoxypyrimidine	410-050-9	5734-64-5	Xn; R22	Xn R: 22 S: (2-)22		

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613-155-00-X	5-chloro-2,3-difluoropyridine	410-090-7	89402-43-7	R10 Xn; R22 R52-53	Xn R: 10-22-52/53 S: (2-)23-36-61		
613-156-00-5	2-butyl-4-chloro-5-formylimidazole	410-260-0	83857-96-9	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
613-157-00-0	2,4-diamino-5-methoxymethylpyrimidine	410-330-0	54236-98-5	Xn; R22-48/22 Xi; R36	Xn R: 22-36-48/22 S: (2-)22-26-36		
613-158-00-6	2,3-dichloro-5-trifluoromethyl-pyridine	410-340-5	69045-84-7	Xn; R20/22 Xi; R41 R43 N; R51-53	Xn; N R: 20/22-41-43-51/53 S: (2-)24-26-37/39-61		
613-159-00-1	fenazaquin (ISO); 4-[2-[4-(1,1-dimethylethyl)phenyl]-ethoxy]quinazoline	410-580-0	120928-09-8	T; R25 Xn; R20 N; R50-53	T; N R: 20-25-50/53 S: (1/2-)37-45-60-61		
613-160-00-7	(1S)-2-methyl-2,5-diazobicyclo[2.2.1]heptane dihydrobromide	411-000-9	125224-62-6	R43	Xi R: 43 S: (2-)24-37		
613-163-00-3	azimsulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-[1-methyl-4-(2-methyl-2*H*-tetrazol-5-yl)pyrazol-5-ylsulfonyl]urea	—	120162-55-2	N; R50-53	N R: 50/53 S: 60-61		
613-164-00-9	flufenacet (ISO); *N*-(4-fluorophenyl)-*N*-isopropyl-2-(5-trifluoromethyl-[1,3,4]thiadiazol-2-yloxy)acetamide	—	142459-58-3	Xn; R22-48/22 R43 N; R50-53	Xn; N R: 22-43-48/22-50/53 S: (2-)13-24-37-60-61		
613-165-00-4	flupyrsulfuron-methyl-sodium (ISO); methyl 2-[[(4,6-dimethoxypyrimidin-2-ylcarbamoyl)sulfamoyl]-6-trifluoromethyl]nicotinate, monosodium salt	—	144740-54-5	N; R50-53	N R: 50/53 S: 60-61		
613-166-00-X	flumioxazin (ISO); *N*-(7-fluoro-3,4-dihydro-3-oxo-4-prop-2-ynyl-2*H*-1,4-benzoxazin-6-yl)cyclohex-1-ene-1,2-dicarboxamide	—	103361-09-7	Repr. Cat. 2; R61 N; R50-53	T; N R: 61-50/53 S: 53-45-60-61		

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613-167-00-5	reaction mass of: 5-chloro-2-methyl-4-isothiazolin-3-one [EC no. 247-500-7] and 2-methyl-2*H* -isothiazol-3-one [EC no. 220-239-6] (3:1); reaction mass of: 5-chloro-2-methyl-4-isothiazolin-3-one [EC no. 247-500-7] and 2-methyl-4-isothiazolin-3-one [EC no. 220-239-6] (3:1)	—	55965-84-9	T; R23/24/25 C; R34 R43 N; R50-53	T; N R: 23/24/25-34-43-50/53 S: (2-)26-28-36/37/39-45-60-61	C; R34: C ≥ 0,6 % Xi; R36/38: 0,06 % ≤ C < 0,6 % R43: C ≥ 0,0015 %	
613-168-00-0	1-vinyl-2-pyrrolidone	201-800-4	88-12-0	Carc. Cat. 3; R40 Xn; R20/21/22-48/20 Xi; R37-41	Xn R: 20/21/22-37-40-41-48/20 S: (2-)26-36/37/39		D
613-169-00-6	9-vinylcarbazole	216-055-0	1484-13-5	Muta. Cat. 3; R68 Xn; R21/22 Xi; R38 R43 N; R50-53	Xn; N R: 21/22-38-43-50/53-68 S: 22-23-36/37-60-61		
613-170-00-1	2,2-ethylmethylthiazolidine	404-500-3	694-64-4	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)24-26-37/39-61		
613-171-00-7	hexaconazole (ISO); (*RS*)-2-(2,4-dichlorophenyl)-1-(1*H*-1,2,4-triazol-1-yl)hexan-2-ol	413-050-7	79983-71-4	Xn; R22 R43 N; R51-53	Xn; N R: 22-43-51/53 S: (2-)24-37-61		
613-172-00-2	5-chloro-1,3-dihydro-2*H*-indol-2-one	412-200-9	17630-75-0	Repr. Cat. 3; R62 Xn; R22 R43 R52-53	Xn R: 22-43-62-52/53 S: (2-)22-36/37-61		
613-173-00-8	fluquinconazole (ISO); 3-(2,4-dichlorophenyl)-6-fluoro-2-(1*H*-1,2,4-triazol-1-yl)quinazolin-4-(3*H*)-one	411-960-9	136426-54-5	T; R23/25-48/25 Xn; R21 Xi; R38 N; R50-53	T; N R: 21-23/25-38-48/25-50/53 S: (1/2-)36/37/39-38-45-60-61		
613-174-00-3	(±) 2-(2,4-dichlorophenyl)-3-(1*H*-1,2,4-triazol-1-yl)propyl-1,1,2,2-tetrafluoroethylether	407-760-6	112281-77-3	Carc. Cat. 3; R40 Xn; R20/22 N; R51-53	Xn; N R: 20/22-40-51/53 S: (2-)36/37-41-61		

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613-175-00-9	epoxiconazole (ISO); (2*RS*,3*SR*)-3-(2-chlorophenyl)-2-(4-fluorophenyl)-[(1*H*-1,2,4-triazol-1-yl)methyl]oxirane	406-850-2	133855-98-8	Carc. Cat. 3; R40 Repr. Cat. 3; R62 Repr. Cat. 3; R63 N; R51-53	Xn; N R: 40-62-63-51/53 S: (2-)36/37-46-61		
613-176-00-4	2-methyl-2-azabicyclo[2.2.1]heptane	404-810-9	4524-95-2	R10 Xn; R21/22-48/20 C; R34	C R: 10-21/22-34-48/20 S: (1/2-)16-26-36/37/39-45		
613-177-00-X	8-amino-7-methylquinoline	412-760-4	5470-82-6	Xn; R21/22 R43 N; R51-53	Xn; N R: 21/22-43-51/53 S: (2-)36/37-61		
613-178-00-5	4-ethyl-2-methyl-2-isopentyl-1,3-oxazolidine	410-470-2	137796-06-6	C; R34 R43	C R: 34-43 S: (1/2-)7/8-26-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
613-179-00-0	lithium 3-oxo-1,2(2*H*)-benzisothiazol-2-ide	411-690-1	111337-53-2	Xn; R22 C; R34 R43 N; R51-53	C; N R: 22-34-43-51/53 S: (1/2-)26-36/37/39-45-61		
613-180-00-6	*N*-(1,1-dimethylethyl)bis(2-benzothiazolesulfen)amide	407-430-1	3741-80-8	N; R50-53	N R: 50/53 S: 60-61		
613-181-00-1	5,5-dimethyl-perhydro-pyrimidin-2-one α-(4-trifluoromethyl-styryl)-α-(4-trifluoromethyl)cinnamylidenehydrazone	405-090-9	67485-29-4	T; R48/25 Xn; R22 Xi; R36 N; R50-53	T; N R: 22-36-48/25-50/53 S: (1/2-)22-26-36/37-45-60-61		
613-182-00-7	1-(1-naphthylmethyl)quinolinium chloride	406-220-7	65322-65-8	Carc. Cat. 3; R40 Muta. Cat. 3; R68 Xn; R22 Xi; R38-41 R52-53	Xn R: 22-38-40-41-52/53-68 S: (2-)22-26-36/37/39-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
613-183-00-2	reaction mass of: 5-(*N*-methylperfluorooctylsulfonamido) methyl-3-octadecyl-1,3-oxazolidin-2-one; 5-(*N*-methylperfluoroheptylsulfonamido)methyl-3-octadecyl-1,3-oxazolidin-2-one	413-640-4	—	Xn; R48/22 N; R50-53	Xn; N R: 48/22-50/53 S: (2-)36-60-61		
613-184-00-8	nitrilotriethyleneammoniopropane-2-ol 2-ethylhexanoate	413-670-8	—	Xi; R36 R43	Xi R: 36-43 S: (2-)24-26-37		
613-185-00-3	2,3,5,6-tetrahydro-2-methyl-2*H*-cyclopenta[*d*]-1,2-thiazol-3-one	407-630-9	82633-79-2	T; R25 Xi; R41 R43 N; R50-53	T; N R: 25-41-43-50/53 S: (1/2-)22-26-36/37/39-45-60-61		
613-186-00-9	(2*R*,3*R*)-3-((*R*)-1-(*tert*-butyldimethylsiloxy)ethyl)-4-oxoazetidin-2-yl acetate	408-050-9	76855-69-1	Xi; R36 R43 N; R51-53	Xi; N R: 36-43-51/53 S: (2-)24-26-37-61		
613-188-00-X	1-(3-(4-fluorophenoxy)propyl)-3-methoxy-4-piperidinone	411-500-7	116256-11-2	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)22-24-26-37/39-61		
613-189-00-5	1,4,7,10-tetrakis(*p*-toluensulfonyl)-1,4,7,10-tetraazacyclododecane	414-030-0	52667-88-6	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
613-190-00-0	disodium 1-amino-4-(2-(5-chloro-6-fluoro-pyrimidin-4-ylamino-methyl)-4-methyl-6-sulfo-phenylamino)-9,10-dioxo-9,10-dihydro-anthracene-2-sulfonate	414-040-5	149530-93-8	Xn; R22 R43	Xn R: 22-43 S: (2-)22-24-37		
613-191-00-6	3-ethyl-2-methyl-2-(3-methylbutyl)-1,3-oxazolidine	421-150-7	143860-04-2	Repr. Cat. 2; R60 C; R34 N; R50-53	T; N R: 60-34-50/53 S: 53-45-60-61		
613-193-00-7	pentakis[3-(dimethylammonio)propylsulfamoyl]-[(6-hydroxy-4,4,8,8-tetramethyl-4,8-diazoniaundecane-1,11-diyldisulfamoyl)di[phthalocyaninecopper(II)]] heptalactate	414-930-3	—	N; R51-53	N R: 51/53 S: 61		
613-194-00-2	6,13-dichloro-3,10-bis{2-[4-fluoro-6-(2-sulfophenylamino)-1,3,5-triazin-2-ylamino]propylamino}benzo[5,6][1,4]oxazino[2,3-.b.]phenoxazine-4,11-disulphonic acid, lithium-, sodium salt	418-000-8	163062-28-0	Xi; R41	Xi R: 41 S: (2-)22-26-39		

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613-195-00-8	2,2-(1,4-phenylene)bis((4*H*-3,1-benzoxazine-4-one)	418-280-1	18600-59-4	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
613-196-00-3	5-[[4-chloro-6-[[2-[[4-fluoro-6-[[5-hydroxy-6-[(4-methoxy-2-sulfophenyl)azo]-7-sulfo-2-naphthalenyl]amino]-1,3,5-triazin-2-yl]amino]-1-methylethyl]amino]-1,3,5-triazin-2-yl]amino]-3-[[4-(ethenylsulfonyl)phenyl]azo]-4-hydroxy-naphtalene-2,7-disulfonic acid, sodium salt	418-380-5	168113-78-8	Xi; R41	Xi R: 41 S: (2-)26-39		
613-197-00-9	reaction mass of: 2,4,6-tri(butylcarbamoyl)-1,3,5-triazine; 2,4,6-tri(methylcarbamoyl)-1,3,5-triazine; [(2-butyl-4,6-dimethyl)tricarbamoyl]-1,3,5-triazine; [(2,4-dibutyl-6-methyl)tricarbamoyl]-1,3,5-triazine	420-390-1	187547-46-2	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
613-199-00-X	reaction mass of: 1,3,5-tris(3-aminomethylphenyl)-1,3,5-(1*H*,3*H*,5*H*)-triazine-2,4,6-trione; reaction mass of oligomers of 3,5-bis(3-aminomethylphenyl)-1-poly[3,5-bis(3-aminomethylphenyl)-2,4,6-trioxo-1,3,5-(1*H*,3*H*,5*H*)-triazin-1-yl]-1,3,5-(1*H*,3*H*,5*H*)-triazine-2,4,6-trione	421-550-1	—	Carc. Cat. 2; R45 Repr. Cat. 2; R61 R43 R52-53	T R: 45-61-43-52/53 S: 53-45-61		
613-200-00-3	Reaction product of: copper, (29*H*,31*H*-phthalocyaninato(2-)-*N*29,*N*30,*N*31,*N*32)-, chlorosulfuric acid and 3-(2-sulfooxyethylsulfonyl)aniline, sodium salts	420-980-7	—	Xi; R41	Xi R: 41 S: (2-)22-26-39		
613-201-00-9	(*R*)-5-bromo-3-(1-methyl-2-pyrrolidinyl methyl)-1*H*-indole	422-390-5	143322-57-0	Repr. Cat. 3; R62 T; R39-48/25 Xn; R20/22 Xi; R41 R43 N; R50-53	T; N R: 20/22-39-41-43-48/25-62-50/53 S: (1/2-)53-45-60-61		
613-202-00-4	pymetrozine (ISO); (*E*)-4,5-dihydro-6-methyl-4-(3-pyridylmethyleneamino)-1,2,4-triazin-3(2*H*)-one	—	123312-89-0	Carc. Cat. 3; R40 R52-53	Xn R: 40-52/53 S: (2-)36/37-61		
613-203-00-X	pyraflufen-ethyl; [1] pyraflufen [2]	[1] [2]	129630-19-9 [1] 129630-17-7 [2]	N; R50-53	N R: 50/53 S: 60-61		

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613-204-00-5	oxadiargyl (ISO); 3-[2,4-dichloro-5-(2-propynyloxy)phenyl]-5-(1,1-dimethylethyl)-1,3,4-oxadiazol-2(3*H*)-one; 5-*tert*-butyl-3-[2,4-dichloro-5-(prop-2-ynyloxy)phenyl]-1,3,4-oxadiazol-2(3*H*)-one	254-637-6	39807-15-3	Repr. Cat. 3; R63 Xn; R48/22 N; R50-53	Xn; N R: 48/22-63-50/53 S: (2-)36/37-46-60-61		
613-205-00-0	propiconazole (ISO); (±)-1-[2-(2,4-dichlorophenyl)-4-propyl-1,3-dioxolan-2-ylmethyl]-1*H*-1,2,4-triazole	262-104-4	60207-90-1	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)36/37-46-60-61		
613-206-00-6	fenamidone (ISO); (*S*)-5-methyl-2-methylthio-5-phenyl-3-phenylamino-3,5-dihydroimidazol-4-one	—	161326-34-7	N; R50-53	N R: 50/53 S: 60-61		
613-208-00-7	imazamox (ISO); (*RS*)-2-(4-isopropyl-4-methyl-5-oxo-2-imidazolin-2-yl)-5-methoxymethylnicotinic acid	—	114311-32-9	N; R50-53	N R: 50/53 S: 60-61		
613-209-00-2	*cis*-1-(3-chloropropyl)-2,6-dimethyl-piperidin hydrochloride	417-430-3	63645-17-0	T; R25 Xn; R48/22 R43 N; R51-53	T; N R: 25-43-48/22-51/53 S: (1/2-)22-36/37-45-61		
613-210-00-8	2-(3-chloropropyl)-2,5,5-trimethyl-1,3-dioxane	417-650-1	88128-57-8	Xn; R48/22 R52-53	Xn R: 48/22-52/53 S: (2-)23-25-36-61		
613-211-00-3	*N*-methyl-4-(*p*-formylstyryl)pyridinium methylsulfate	418-240-3	74401-04-0	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
613-212-00-9	4-[4-(2-ethylhexyloxy)phenyl](1,4-thiazinane-1,1-dioxide)	418-320-8	133467-41-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-60-61		
613-213-00-4	*cis*-1-benzoyl-4-[(4-methylsulfonyl)oxy]-L-proline	416-040-0	120807-02-5	R52-53	R: 52/53 S: 61		
613-214-00-X	*N*,*N*-di-*n*-butyl-2-(1,2-dihydro-3-hydroxy-6-isopropyl-2-quinolylidene)-1,3-dioxoindan-5-carboxamide	416-260-7	147613-95-4	R53	R: 53 S: 61		

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613-215-00-5	2-chloromethyl-3,4-dimethoxypyridinium chloride	416-440-5	72830-09-2	Xn; R21/22-48/22 Xi; R38-41 R43 N; R51-53	Xn; N R: 21/22-38-41-43-48/22-51/53 S: (2-)26-36/37/39-61		
613-216-00-0	6-*tert*-butyl-7-(6-diethylamino-2-methyl-3-pyridylimino)-3-(3-methylphenyl)pyrazolo[3,2-c][1,2,4]triazole	416-490-8	162208-01-7	N; R50-53	N R: 50/53 S: 60-61		
613-217-00-6	4-[3-(3,5-di-*tert*-butyl-4-hydroxyphenyl)propionyloxy]-1-[2-[3-(3,5-di-*tert*-butyl-4-hydrophenyl)propionyloxy]ethyl]-2,2,6,6-tetramethylpiperidine	416-770-1	73754-27-5	R53	R: 53 S: 61		
613-218-00-1	6-hydroxyindole	417-020-4	2380-86-1	Xn; R22 Xi; R41 R43 N; R51-53	Xn; N R: 22-41-43-51/53 S: (2-)24-26-37/39-61		
613-219-00-7	7a-ethyl-3,5-bis(1-methylethyl)-2,3,4,5-tetrahydrooxazolo[3,4-c]-2,3,4,5-tetrahydrooxazole	417-140-7	79185-77-6	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)37-61		
613-220-00-2	trans-(4*S*,6*S*)-5,6-dihydro-6-methyl-4*H*-thieno[2,3-b]thiopyran-4-ol, 7,7-dioxide	417-290-3	147086-81-5	Xn; R22	Xn R: 22 S: (2-)36		
613-221-00-8	2-chloro-5-methyl-pyridine	418-050-0	18368-64-4	Xn; R21/22 Xi; R38 R52-53	Xn R: 21/22-38-52/53 S: (2-)23-25-36/37-61		
613-222-00-3	4-(1-oxo-2-propenyl)-morpholine	418-140-1	5117-12-4	Xn; R22-48/22 Xi; R41 R43	Xn R: 22-41-43-48/22 S: (2-)23-26-36/37/39		
613-223-00-9	*N*-isopropyl-3-(4-fluorophenyl)-1*H*-indole	418-790-4	93957-49-4	R53	R: 53 S: 61		
613-224-00-4	2,5-dimercaptomethyl-1,4-dithiane	419-770-8	136122-15-1	Xn; R22 C; R34 R43 N; R50-53	C; N R: 22-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
613-225-00-X	reaction mass of:[2-(anthraquinon-1-ylamino)-6-[(5-benzoylamino)-anthraquinone-1-ylamino]-4-phenyl]-1,3,5-triazine; 2,6-bis-[(5-benzoylamino)-anthraquinon-1-ylamino]-4-phenyl-1,3,5-triazine.	421-290-9	—	Xn; R48/22 R53	Xn R: 48/22-53 S: (2-)22-36-61		
613-226-00-5	1-(2-(ethyl(4-(4-(4-(4-(ethyl(2-pyridinoethyl)amino)-2-methylphenylazo)benzoylamino)-phenylazo)-3-methylphenyl) amino)ethyl)-pyridinium dichloride	420-950-3	163831-67-2	Xi; R41 N; R50-53	Xi; N R: 41-50/53 S: (2-)26-39-60-61		
613-227-00-0	(±)-[(*R**,*R**) and (*R**,*S**)]-6-fluoro-3,4-dihydro-2-oxiranyl-2*H*-1-benzopyran	419-600-2	99199-90-3	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-28-36/37-61		
613-228-00-6	(±)-(*R**,*S**)-6-fluoro-3,4-dihydro-2-oxiranyl-2*H*-1-benzopyran	419-630-6	793669-26-8	N; R51-53	N R: 51/53 S: 24-61		
613-230-00-7	florasulam (ISO); 2',6',8-trifluoro-5-methoxy-5-triazolo[1,5-c]; pyrimidine-2-sulfonanilide	—	145701-23-1	N; R50-53	N R: 50/53 S: 60-61		
613-233-00-3	4,4'-(oxy-(bismethylene))-bis-1,3-dioxolane	423-230-7	56552-15-9	Xi; R41	Xi R: 41 S: (2-)26-39		
614-001-00-4	nicotine (ISO); 3-(*N*-methyl-2-pyrrolidinyl)pyridine	200-193-3	54-11-5	T+; R27 T; R25 N; R51-53	T+; N R: 25-27-51/53 S: (1/2-)36/37-45-61		
614-002-00-X	salts of nicotine	—	—	T+; R26/27/28 N; R51-53	T+; N R: 26/27/28-51/53 S: (1/2-)13-28-45-61		A
614-003-00-5	strychnine	200-319-7	57-24-9	T+; R27/28 N; R50-53	T+; N R: 27/28-50/53 S: (1/2-)36/37-45-60-61		
614-004-00-0	salts of strychnine	—	—	T+; R26/28 N; R50-53	T+; N R: 26/28-50/53 S: (1/2-)13-28-45-60-61		A

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614-005-00-6	colchicine	200-598-5	64-86-8	T+; R26/28	T+ R: 26/28 S: (1/2-)13-45		
614-006-00-1	brucine; 2,3-dimethoxystrychnine	206-614-7	357-57-3	T+; R26/28 R52-53	T+ R: 26/28-52/53 S: (1/2-)13-45-61		
614-007-00-7	brucine sulphate; [1] brucine nitrate; [2] Strychnidin-10-one, 2,3-dimethoxy-, mono[(R)-1-methylheptyl 1,2-benzenedicarboxylate]; [3] Strychnidin-10-one, 2,3-dimethoxy-, compd. with (*S*)mono(1-methylheptyl)-1,2-benzenedicarboxylate (1:1) [4]	225-432-9 [1] 227-317-9 [2] 269-439-5 [3] 269-710-8 [4]	4845-99-2 [1] 5786-97-0 [2] 68239-26-9 [3] 68310-42-9 [4]	T+; R26/28 R52-53	T+ R: 26/28-52/53 S: (1/2-)13-45-61		A
614-008-00-2	aconitine	206-121-7	302-27-2	T+; R26/28	T+ R: 26/28 S: (1/2-)24-45		
614-009-00-8	salts of aconitine	—	—	T+; R26/28	T+ R: 26/28 S: (1/2-)24-45		A
614-010-00-3	atropine	200-104-8	51-55-8	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		
614-011-00-9	salts of atropine	—	—	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		A
614-012-00-4	hyoscyamine	202-933-0	101-31-5	T+; R26/28	T+ R: 26/28 S: (1/2-)24-45		
614-013-00-X	salts of hyoscyamine	—	—	T+; R26/28	T+ R: 26/28 S: (1/2-)24-45		A
614-014-00-5	hyoscine	200-090-3	51-34-3	T+; R26/27/28	T+ R: 26/27/28 S: (1/2-)25-45		

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614-015-00-0	salts of hyoscine	—	—	T+; R26/27/28	T+ R: 26/27/28 S: (1/2-)25-45		A
614-016-00-6	pilocarpine	202-128-4	92-13-7	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		
614-017-00-1	salts of pilocarpine	—	—	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		A
614-018-00-7	papaverine	200-397-2	58-74-2	Xn; R22	Xn R: 22 S: (2-)22		
614-019-00-2	salts of papaverine	—	—	Xn; R22	Xn R: 22 S: (2-)22		A
614-020-00-8	physostigmine	200-332-8	57-47-6	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		
614-021-00-3	salts of physostigmine	—	—	T+; R26/28	T+ R: 26/28 S: (1/2-)25-45		A
614-022-00-9	digitoxin	200-760-5	71-63-6	T; R23/25 R33	T R: 23/25-33 S: (1/2-)45		
614-023-00-4	ephedrine	206-080-5	299-42-3	Xn; R22	Xn R: 22 S: (2-)22-25		
614-024-00-X	salts of ephedrine	—	—	Xn; R22	Xn R: 22 S: (2-)22-25		A
614-025-00-5	ouabain	211-139-3	630-60-4	T; R23/25 R33	T R: 23/25-33 S: (1/2-)45		

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614-026-00-0	strophantin-K	234-239-9	11005-63-3	T; R23/25 R33	T R: 23/25-33 S: (1/2-)45		
614-027-00-6	bufa-4,20,22-trienolide, 6-(acetyloxy)-3-(β-D-glucopyranosyloxy)-8,14-dihydroxy-, (3β, 6β)-; red squill; scilliroside	208-077-4	507-60-8	T+; R28	T+ R: 28 S: (1/2-)36/37-45		
614-028-00-1	reaction mass of: 2-ethylhexyl mono-D-glucopyranoside; 2-ethylhexyl di-D-glucopyranoside	414-420-0	—	Xi; R41	Xi R: 41 S: (2-)26-39		
614-029-00-7	constitutional isomers of penta-O-allyl-β-D-fructofuranosyl-α-D-glucopyranoside; constitutional isomers of hexa-O-allyl-β-D-fructofuranosyl-α-D-glucopyranoside; constitutional isomers of hepta-O-allyl-β-D-fructofuransoyl-α-D-glucopyranoside	419-640-0	68784-14-5	Xn; R22	Xn R: 22 S: (2-)		
615-001-00-7	methyl isocyanate	210-866-3	624-83-9	F+; R12 ⊗ Repr. Cat. 3; R63 T+; R26 T; R24/25 R42/43 Xi; R37/38-41	F+; T+ R: 12-24/25-26-37/38-41-42/43-63 S: (1/2-)26-27/28-36/37/39-45-63		
615-002-00-2	methyl isothiocyanate	209-132-5	556-61-6	T; R23/25 C; R34 R43 N; R50-53	T; N R: 23/25-34-43-50/53 S: (1/2-)36/37-38-45-60-61		
615-003-00-8	thiocyanic acid	207-337-4	463-56-9	Xn; R20/21/22 R32 R52-53	Xn R: 20/21/22-32-52/53 S: (2-)13-61		
615-004-00-3	salts of thiocyanic acid	—	—	Xn; R20/21/22 R32 R52-53	Xn R: 20/21/22-32-52/53 S: (2-)13-61		A

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615-005-00-9	4,4'-methylenediphenyl diisocyanate; diphenylmethane-4,4'-diisocyanate; [1] 2,2'-methylenediphenyl diisocyanate; diphenylmethane-2,2'-diisocyanate; [2] *o*-(*p*-isocyanatobenzyl)phenyl isocyanate; diphenylmethane-2,4'-diisocyanate; [3] methylenediphenyl diisocyanate [4]	202-966-0 [1] 219-799-4 [2] 227-534-9 [3] 247-714-0 [4]	101-68-8 [1] 2536-05-2 [2] 5873-54-1 [3] 26447-40-5 [4]	Xn; R20 Xi; R36/37/38 R42/43	Xn R: 20-36/37/38-42/43 S: (1/2-)23-36/37-45	Xi; R36/37/38: C ≥ 5 % R42: C ≥ 0,1 %	C 2
615-006-00-4	2-methyl-*m*-phenylene diisocyanate; toluene-2,4-di-isocyanate; [1] 4-methyl-*m*-phenylene diisocyanate; toluene-2,6-di-isocyanate; [2] *m*-tolylidene diisocyanate; toluene-diisocyanate [3]	202-039-0 [1] 209-544-5 [2] 247-722-4 [3]	91-08-7 [1] 584-84-9 [2] 26471-62-5 [3]	Carc. Cat. 3; R40 T+; R26 Xi; R36/37/38 R42/43 R52-53	T+ R: 26-36/37/38-40-42/43-52/53 S: (1/2-)23-36/37-45-61	R42: C ≥ 0,1 %	C 2
615-007-00-X	1,5-naphthylene diisocyanate	221-641-4	3173-72-6	Xn; R20 Xi; R36/37/38 R42 R52-53	Xn R: 20-36/37/38-42-52/53 S: (2-)26-28-38-45-61		
615-008-00-5	3-isocyanatomethyl-3,5,5-trimethylcyclohexyl isocyanate; isophorone di-isocyanate	223-861-6	4098-71-9	T; R23 Xi; R36/37/38 R42/43 N; R51-53	T; N R: 23-36/37/38-42/43-51/53 S: (1/2-)26-28-38-45-61	T; R23: C ≥ 2 % Xn; R20: 0,5 % ≤ C < 2 % R42/43: C ≥ 0,5 %	2
615-009-00-0	4,4'-methylenedi(cyclohexyl isocyanate); dicyclohexylmethane-4,4'-di-isocyanate	225-863-2	5124-30-1	T; R23 Xi; R36/37/38 R42/43	T R: 23-36/37/38-42/43 S: (1/2-)26-28-38-45	T; R23: C ≥ 2 % Xn; R20: 0,5 % ≤ C < 2 % R42/43: C ≥ 0,5 %	2
615-010-00-6	2,2,4-trimethylhexamethylene-1,6-di-isocyanate; [1] 2,4,4-trimethylhexamethylene-1,6-di-isocyanate [2]	241-001-8 [1] 239-714-4 [2]	16938-22-0 [1] 15646-96-5 [2]	T; R23 Xi; R36/37/38 R42	T R: 23-36/37/38-42 S: (1/2-)26-28-38-45	T; R23: C ≥ 2 % Xn; R20: 0,5 % ≤ C < 2 % R42: C ≥ 0,5 %	C 2
615-011-00-1	hexamethylene-di-isocyanate	212-485-8	822-06-0	T; R23 Xi; R36/37/38 R42/43	T R: 23-36/37/38-42/43 S: (1/2-)26-28-38-45	T; R23: C ≥ 2 % Xn; R20: 0,5 % ≤ C < 2 % R42/43: C ≥ 0,5 %	2

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615-012-00-7	4-isocyanatosulphonyltoluene; tosyl isocyanate	223-810-8	4083-64-1	R14 Xi; R36/37/38 R42	Xn R: 14-36/37/38-42 S: (2-)26-28-30	Xi; R36/37/38: C ≥ 5 %	
615-013-00-2	cyanamide; carbanonitril	206-992-3	420-04-2	T; R25 Xn; R21 Xi; R36/38 R43	T R: 21-25-36/38-43 S: (1/2-)3-22-36/37-45		
615-014-00-8	tris(1-dodecyl-3-methyl-2-phenylbenzimidazolium)hexacyanoferrate	—	7276-58-6	Xn; R22	Xn R: 22 S: (2-)24		
615-015-00-3	1,7,7-trimethylbicyclo(2,2,1)hept-2-yl thiocyanatoacetate; isobornyl thiocyanoacetate	204-081-5	115-31-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)24/25-60-61		
615-016-00-9	potassium cyanate	209-676-3	590-28-3	Xn; R22	Xn R: 22 S: (2-)24/25		
615-017-00-4	calcium cyanamide	205-861-8	156-62-7	Xn; R22 Xi; R37-41	Xn R: 22-37-41 S: (2-)22-26-36/37/39		
615-018-00-X	2-(2-butoxyethoxy)ethyl thiocyanate	203-985-7	112-56-1	R10 T; R24/25	T R: 10-24/25 S: (1/2-)13-36/37-45		
615-019-00-5	dicyclohexylcarbodiimide	208-704-1	538-75-0	T; R24 Xn; R22 Xi; R41 R43	T R: 22-24-41-43 S: (1/2-)24-26-37/39-45		
615-020-00-0	methylene dithiocyanate	228-652-3	6317-18-6	T+; R26 T; R25 C; R34 R43 N; R50	T+; N R: 25-26-34-43-50 S: (1/2-)26-28-36/37/39-45-61		

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615-021-00-6	1,3,5-tris(oxiranylmethyl)-1,3,5-triazine-2,4,6(1*H*,3*H*,5*H*)-trione; TGIC	219-514-3	2451-62-9	Muta. Cat. 2; R46 T; R23/25 Xn; R48/22 Xi; R41 R43 R52-53	T R: 46-23/25-41-43-48/22-52/53 S: 53-45-61		E
615-022-00-1	methyl 3-isocyanatosulfonyl-2-thiophene-carboxylate	410-550-7	79277-18-2	E; R2 ⊗ R14 Xn; R48/22 R42/43	E; Xn R: 2-14-42/43-48/22 S: (2-)22-30-35-36/37		
615-023-00-7	2-(isocyanatosulfonylmethyl)benzoic acid methyl ester; (alt.):methyl 2-(isocyanatosulfonylmethyl)benzoate	410-900-9	83056-32-0	R10 R14 Muta. Cat. 3; R68 Xn; R20-48/22 Xi; R41 R42	Xn R: 10-14-20-41-42-48/22-68 S: (2-)23-26-36/37/39		
615-024-00-2	2-phenylethylisocyanate	413-080-0	1943-82-4	T; R23 Xn; R22 C; R35 R42/43 N; R51-53	T; C; N R: 22-23-35-42/43-51/53 S: (1/2-)23-26-36/37/39-43-45-61		
615-025-00-8	4,4'-ethylidenediphenyl dicyanate	405-740-1	47073-92-7	Xn; R20/22-48/22 Xi; R41 N; R50-53	Xn; N R: 20/22-41-48/22-50/53 S: (2-)26-36/37/39-60-61		
615-026-00-3	4,4'-methylenebis(2,6-dimethylphenyl cyanate)	405-790-4	101657-77-6	R43 R52-53	Xi R: 43-52/53 S: (2-)22-24-37-61		
615-028-00-4	ethyl 2-(isocyanatosulfonyl)benzoate	410-220-2	77375-79-2	E; R2 ⊗ R14 Xn; R22-48/22 Xi; R41 R42/43	E; Xn R: 2-14-22-41-42/43-48/22 S: (2-)8-23-26-30-35-36/37/39		

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615-029-00-X	2,5-bis-isocyanatomethyl-bicyclo[2.2.1]heptane	411-280-2	—	T+; R26 Xn; R22 C; R34 R42/43 R52-53	T+ R: 22-26-34-42/43-52/53 S: (1/2-)23-26-28-36/37/39-45-61		
615-030-00-5	alkali salts, alkali earth salts and other salts of thiocyanic acid not mentioned elsewhere in this Annex	—	—	Xn; R20/21/22 R32 R52-53	Xn R: 20/21/22-32-52/53 S: (2-)13-61		A
615-031-00-0	thallium salt of thiocyanic acid	222-571-7	3535-84-0	Xn; R20/21/22 R32 N; R51-53	Xn; N R: 20/21/22-32-51/53 S: (2-)13-61		A
615-032-00-6	metal salts of thiocyanic acid not mentioned elsewhere in this Annex	—	—	Xn; R20/21/22 R32 N; R50-53	Xn; N R: 20/21/22-32-50/53 S: (2-)13-60-61		A
616-001-00-X	*N*,*N*-dimethylformamide; dimethyl formamide	200-679-5	68-12-2	Repr. Cat. 2; R61 Xn; R20/21 Xi; R36	T R: 61-20/21-36 S: 53-45		E
616-002-00-5	2-fluoroacetamide	211-363-1	640-19-7	T+; R28 T; R24	T+ R: 24-28 S: (1/2-)36/37-45		
616-003-00-0	acrylamide; prop-2-enamide	201-173-7	79-06-1	Carc. Cat. 2; R45 Muta. Cat. 2; R46 Repr. Cat. 3; R62 T; R25-48/23/24/25 Xn; R20/21 Xi; R36/38 R43	T R: 45-46-20/21-25-36/38-43-48/23/24/25-62 S: 53-45		DE

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616-004-00-6	allidochlor (ISO); *N*,*N*-diallylchloroacetamide	202-270-7	93-71-0	Xn; R21/22 Xi; R36/38 N; R51-53	Xn; N R: 21/22-36/38-51/53 S: (2-)26-28-36/37/39-61		
616-005-00-1	chlorthiamid (ISO); 2,6-dichloro (thiobenzamide)	217-637-7	1918-13-4	Xn; R22	Xn R: 22 S: (2-)36		
616-006-00-7	dichlofluanid (ISO); *N*-dichlorofluoromethylthio-*N*',*N*'-dimethyl-*N*-phenylsulphamide	214-118-7	1085-98-9	Xn; R20 Xi; R36 R43 N; R50-53	Xn; N R: 20-36-43-50/53 S: (2-)24-37-60-61		
616-007-00-2	diphenamid (ISO); *N*,*N*-dimethyl-2,2-diphenylacetamide	213-482-4	957-51-7	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
616-008-00-8	propachlor (ISO); 2-chloro-*N*-isopropylacetanilide; α-chloro-*N*-isopropylacetanilide	217-638-2	1918-16-7	Xn; R22 Xi; R36 R43 N; R50-53	Xn; N R: 22-36-43-50/53 S: (2-)24-37-60-61		
616-009-00-3	propanil (ISO); 3',4'-dichloropropionanilide	211-914-6	709-98-8	Xn; R22 N; R50	Xn; N R: 22-50 S: (2-)22-61		
616-010-00-9	tosylchloramide sodium	204-854-7	127-65-1	Xn; R22 R31 C; R34 R42	C R: 22-31-34-42 S: (1/2-)7-22-26-36/37/39-45		
616-011-00-4	*N*,*N*-dimethylacetamide	204-826-4	127-19-5	Repr. Cat. 2; R61 Xn; R20/21	T R: 61-20/21 S: 53-45	Repr. Cat. 2; R61: C ≥ 5 %	E
616-012-00-X	*N*-(dichlorofluoromethylthio)phthalimide; *N*-(fluorodichloromethylthio)phthalimide	211-952-3	719-96-0	Xi; R38	Xi R: 38 S: (2-)28		
616-013-00-5	butyraldehyde oxime	203-792-8	110-69-0	T; R24 Xn; R22 Xi; R36	T R: 22-24-36 S: (1/2-)23-36-45		

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616-014-00-0	2-butanone oxime; ethyl methyl ketoxime; ethyl methyl ketone oxime	202-496-6	96-29-7	Carc. Cat. 3; R40 Xn; R21 Xi; R41 R43	Xn R: 21-40-41-43 S: (2-)13-23-26-36/37/39		
616-015-00-6	alachlor (ISO); 2-chloro-2',6'-diethyl-*N*-(methoxymethyl)acetanilide	240-110-8	15972-60-8	Carc. Cat. 3; R40 Xn; R22 R43 N; R50-53	Xn; N R: 22-40-43-50/53 S: (2-)36/37-46-60-61	N; R50-53: C ≥ 0,2,5 % N; R51-53: 0,025 % ≤ C < 0,25 % R52-53: 0,0025 % ≤ C < 0,025 %	
616-016-00-1	1-(3,4-dichlorophenylimino) thiosemicarbazide	—	5836-73-7	T+; R28	T+ R: 28 S: (1/2-)22-36/37-45		
616-017-00-7	cartap hydrochloride	239-309-2	15263-52-2	Xn; R21/22 N; R50-53	Xn; N R: 21/22-50/53 S: (2-)36/37-60-61		
616-018-00-2	*N*,*N*-diethyl-*m*-toluamide; deet	205-149-7	134-62-3	Xn; R22 Xi; R36/38 R52-53	Xn R: 22-36/38-52/53 S: (2-)61		
616-019-00-8	perfluidone (ISO); 1,1,1-trifluoro-*N*-(4-phenylsulphonyl-*o*-tolyl)methanesulphonamide	253-718-3	37924-13-3	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)		
616-020-00-3	tebuthiuron (ISO); 1-(5-*tert*-butyl-1,3,4-thiadiazol-2-yl)-1,3-dimethylurea	251-793-7	34014-18-1	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)37-60-61		
616-021-00-9	thiazafluron (ISO); 1,3-dimethyl-1-(5-trifluoromethyl-1,3,4-thiadiazol-2-yl)urea	246-901-4	25366-23-8	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)60-61		
616-022-00-4	acetamide	200-473-5	60-35-5	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		
616-023-00-X	*N*-hexadecyl(or octadecyl)-*N*-hexadecyl(or octadecyl)benzamide	401-980-6	—	Xi; R38 R43	Xi R: 38-43 S: (2-)24-37		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
616-024-00-5	2-(4,4-dimethyl-2,5-dioxooxazolidin-1-yl)-2-chloro-5-(2-(2,4-di-*tert*-pentylphenoxy)butyramido)-4,4-dimethyl-3-oxovaleranilide	402-260-4	54942-74-4	R53	R: 53 S: 61		
616-025-00-0	valinamide	402-840-7	20108-78-5	Repr. Cat. 3; R62 Xi; R36 R43	Xn R: 36-43-62 S: (2-)26-36/37		
616-026-00-6	thioacetamide	200-541-4	62-55-5	Carc. Cat. 2; R45 Xn; R22 Xi; R36/38 R52-53	T R: 45-22-36/38-52/53 S: 53-45-61		E
616-027-00-1	tris(2-(2-hydroxyethoxy)ethyl)ammonium 3-acetoacetamido-4-methoxybenzenesulfonate	403-760-5	—	R43	Xi R: 43 S: (2-)24-37		
616-028-00-7	*N*-(4-(3-(4-cyanophenyl)ureido)-3-hydroxyphenyl)-2-(2,4-di-*tert*-pentylphenoxy)octanamide	403-790-9	108673-51-4	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
616-029-00-2	*N*,*N*'-ethylenebis(vinylsulfonylacetamide)	404-790-1	66710-66-5	Xi; R41 R43	Xi R: 41-43 S: (2-)24-26-37/39		
616-030-00-8	ethidimuron (ISO); 1-(5-ethylsulphonyl-1,3,4-thiadiazol-2-yl)-1,3-dimethylurea	250-010-6	30043-49-3	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
616-031-00-3	dimethachlor (ISO); 2-chloro-*N*-(2,6-dimethylphenyl)-*N*-(2-methoxyethyl)acetamide	256-625-6	50563-36-5	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)24-37-60-61		
616-032-00-9	diflufenican (ISO); *N*-(2,4-difluorophenyl)-2-[3-(trifluoromethyl)phenoxy]-3-pyridinecarboxamide	—	83164-33-4	R52-53	R: 52/53 S: 61		
616-033-00-4	cyprofuram (ISO); *N*-(3-chlorophenyl)-*N*-(tetrahydro-2-oxo-3-furyl)cyclopropanecarboxamide	274-050-9	69581-33-5	T; R25 Xn; R21 N; R50-53	T; N R: 21-25-50/53 S: (1/2-)36/37-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
616-034-00-X	pyracarbolid (ISO); 3,4-dihydro-6-methyl-2*H*-pyran-5-carboxanilide	246-419-4	24691-76-7	R52-53	R: 52/53 S: 61		
616-035-00-5	cymoxanil (ISO); 2-cyano-*N*-[(ethylamino)carbonyl]-2-(methoxyimino)acetamide	261-043-0	57966-95-7	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)36/37-60-61		
616-036-00-0	2-chloracetamide	201-174-2	79-07-2	Repr. Cat. 3; R62 T; R25 R43	T R: 25-43-62 S: (1/2-)22-36/37-45	R43: C ≥ 0,1 %	
616-037-00-6	acetochlor (ISO); 2-chloro-*N*-(ethoxymethyl)-*N*-(2-ethyl-6-methylphenyl)acetamide	251-899-3	34256-82-1	Xn; R20 Xi; R37/38 R43 N; R50-53	Xn; N R: 20-37/38-43-50/53 S: (2-)36/37-60-61		
616-038-00-1	(4-aminophenyl)-*N*-methylmethylensulfonamide hydrochloride	406-010-5	88918-84-7	Xi; R41 R43 N; R51-53	Xi; N R: 41-43-51/53 S: (2-)24-26-37/39-61		
616-039-00-7	3',5'-dichloro-4'-ethyl-2'-hydroxypalmitanilide	406-200-8	117827-06-2	R43	Xi R: 43 S: (2-)24-37		
616-040-00-2	potassium *N*-(4-toluenesulfonyl)-4-toluenesulfonamide	406-650-5	97888-41-0	Xi; R41	Xi R: 41 S: (2-)26-39		
616-041-00-8	3',5'-dichloro-2-(2,4-di-*tert*-pentylphenoxy)-4'-ethyl-2'-hydroxyhexananilide	406-840-8	101664-25-9	R53	R: 53 S: 61		
616-042-00-3	*N*-(2-(6-ethyl-7-(4-methylphenoxy)-1*H*-pyrazolo[1,5-b][1,2,4]triazol-2-yl)propyl)-2-octadecyloxybenzamide	407-070-5	142859-67-4	R43 R53	Xi R: 43-53 S: (2-)22-24-37-61		
616-043-00-9	isoxaben (ISO); *N*-[3-(1-ethyl-1-methylpropyl)-1,2-oxazol-5-yl]-2,6-dimethoxybenzamide	407-190-8	82558-50-7	R53	R: 53 S: 61		
616-044-00-4	*N*-(3,5-dichloro-4-ethyl-2-hydroxyphenyl)-2-(3-pentadecylphenoxy)-butanamide	402-510-2	—	N; R51-53	N R: 51/53 S: 61		

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616-045-00-X	2'-(4-chloro-3-cyano-5-formyl-2-thienylazo)-5'-diethylamino-2-methoxyacetanilide	405-190-2	122371-93-1	R43 R53	Xi R: 43-53 S: 2-22-24-37-61		
616-046-00-5	*N*-(2-(6-chloro-7-methylpyrazolo(1,5-b)-1,2,4-triazol-4-yl) propyl)-2-(2,4-di-*tert*-pentylphenoxy)octanamide	406-390-2	—	N; R50-53	N R: 50/53 S: 60-61		
616-047-00-0	reaction mass of: 2,2',2",2"'-(ethylenedinitrilotetrakis-*N*,*N*-di (C_{16})alkylacetamide; 2,2',2",2"'-(ethylenedinitrilotetrakis-*N*,*N*-di(C_{18})alkylacetamide	406-640-0	—	R43	Xi R: 43 S: (2-)24-37		
616-048-00-6	3'-trifluoromethylisobutyranilide	406-740-4	1939-27-1	Xn; R48/22 N; R51-53	Xn; N R: 48/22-51/53 S: (2-)22-36-61		
616-049-00-1	2-(2,4-bis(1,1-dimethylethyl)phenoxy)-*N*-(3,5-dichloro-4-ethyl-2-hydroxyphenyl)-hexanamide	408-150-2	99141-89-6	R53	R: 53 S: 61		
616-050-00-7	lufenuron (ISO); *N*-[2,5-dichloro-4-(1,1,2,3,3,3-hexafluoropropoxy)-phenyl-aminocarbonyl]-2,6-difluorobenzamide	410-690-9	103055-07-8	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
616-051-00-2	reaction mass of: 2,4 -bis(*N*'-(4-methylphenyl)-ureido)-toluene; 2,6 -bis(*N*'-(4-methylphenyl)-ureido)-toluene	411-070-0	—	R53	R: 53 S: 61		
616-052-00-8	formamide	200-842-0	75-12-7	Repr. Cat. 2; R61	T R: 61 S: 53-45		
616-053-00-3	*N*-methylacetamide	201-182-6	79-16-3	Repr. Cat. 2; R61	T R: 61 S: 53-45		
616-054-00-9	iprodione (ISO); 3-(3,5-dichlorophenyl)-2,4-dioxo-*N*-isopropylimidazolidine-1-carboxamide	253-178-9	36734-19-7	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		
616-055-00-4	propyzamide (ISO); 3,5-dichloro-*N*-(1,1-dimethylprop-2-ynyl)benzamide	245-951-4	23950-58-5	Carc. Cat. 3; R40 N; R50-53	Xn; N R: 40-50/53 S: (2-)36/37-60-61		

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616-056-00-X	*N*-methylformamide	204-624-6	123-39-7	Repr. Cat. 2; R61 Xn; R21	T R: 61-21 S: 53-45		E
616-057-00-5	reaction mass of: *N*-[3-hydroxy-2-(2-methylacryloylamino-methoxy)propoxymethyl]-2-methylacrylamide; *N*-[2,3-bis-(2-methylacryloylaminomethoxy)propoxymethyl]-2-methylacrylamide; methacrylamide; 2-methyl-*N*-(2-methylacryloylaminomethoxymethyl)-acrylamide; *N*-(2,3-dihydroxypropoxymethyl)-2-methylacrylamide	412-790-8	—	Carc. Cat. 2; R45 Muta. Cat. 3; R68 Xn; R48/22	T R: 45-48/22 S: 53-45		E
616-058-00-0	1,3-bis(3-methyl-2,5-dioxo-1*H*-pyrrolinylmethyl)benzene	412-570-1	119462-56-5	Xn; R48/22 Xi; R41 R43 N; R50-53	Xn; N R: 41-43-48/22-50/53 S: (2-)26-36/37/39-60-61		
616-059-00-6	4-((4-(diethylamino)-2-ethoxyphenyl)imino)-1,4-dihydro-1-oxo-*N*-propyl-2-naphthalenecarboxamide	412-650-6	121487-83-0	R53	R: 53 S: 61		
616-060-00-1	Condensation product of: 3-(7-carboxyhept-1-yl)-6-hexyl-4-cyclohexene-1,2-dicarboxylic acid with polyamines (primarily amino-ethyl-piperazine and triethylenetetramine)	413-770-1	—	Xn; R22 C; R34 R43 N; R50-53	C; N R: 22-34-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
616-061-00-7	*N*,*N*'-1,6-hexanediylbis(*N*-(2,2,6,6-tetramethyl-piperidin-4-yl)-formamide	413-610-0	124172-53-8	Xi; R36 R52-53	Xi R: 36-52/53 S: (2-)26-61		
616-062-00-2	*N*-[3-[(2-acetyloxy)ethyl](phenyl-methyl)amino]-4-methoxy-phenylacetamide	411-590-8	70693-57-1	C; R34 R52-53	C R: 34-52/53 S: (1/2-)26-36/37/39-45-61		
616-063-00-8	3-dodecyl-(1-(1,2,2,6,6-pentamethyl-4-piperidin)-yl)-2,5-pyrrolidindione	411-920-0	106917-30-0	T; R23 Xn; R22-48/22 C; R35 N; R50-53	T; C; N R: 22-23-35-48/22-50/53 S: (1/2-)26-28-36/37/39-45-60-61		
616-064-00-3	*N*-*tert*-butyl-3-methylpicolinamide	406-720-5	32998-95-1	R52-53	R: 52/53 S: 61		

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616-065-00-9	3'-(3-acetyl-4-hydroxyphenyl)-1,1-diethylurea	411-970-3	79881-89-3	Xn; R22-48/22	Xn R: 22-48/22 S: (2-)22-36		
616-066-00-4	5,6,12,13-tetrachloroanthra(2,1,9-*def*:6,5,10-*d'e'f'*)diisoquinoline-1,3,8,10(2*H*,9*H*)-tetrone	405-100-1	115662-06-1	Repr. Cat. 3; R62	Xn R: 62 S: (2-)22-36/37		
616-067-00-X	dodecyl 3-(2-(3-benzyl-4-ethoxy-2,5-dioxoimidazolidin-1-yl)-4,4-dimethyl-3-oxovaleramido)-4-chlorobenzoate	407-300-4	92683-20-0	R53	R: 53 S: 61		
616-068-00-5	potassium 4-(11-methacrylamidoundecanamido)benzenesulfonate	406-500-9	174393-75-0	R43	Xi R: 43 S: (2-)22-24-37		
616-069-00-0	1-hydroxy-5-(2-methylpropyloxycarbonylamino)-*N*-(3-dodecyloxypropyl)-2-naphthoamide	406-210-2	110560-22-0	R53	R: 53 S: 61		
616-070-00-6	reaction mass of: 3,3'-dicyclohexyl-1,1'-methylenebis(4,1-phenylene)diurea; 3-cyclohexyl-1-(4-(4-(3-octadecylureido)benzyl)phenyl)urea; 3,3'-dioctadecyl-1,1'-methylenebis(4,1-phenylene)diurea	406-530-2	—	R53	R: 53 S: 22-61		

616-071-00-1	reaction mass of: bis(*N*-cyclohexyl-*N'*-phenyleneureido) methylene;
	bis(*N*-octadecyl-*N'*-phenyleneureido)methylene;
	bis(*N*-dicyclohexyl-*N'*-phenyleneureido)methylene (1:2:1)
		406-550-1	—	R43
				R53
					Xi
					R: 43-53
					S: (2-)22-24-37-61		

616-072-00-7	1-(2-deoxy-5-*O*-trityl-β-D-threopentofuranosyl)thymine	407-120-6	55612-11-8	R53	R: 53 S: 61		
616-073-00-2	4'-ethoxy-2-benzimidazoleanilide	407-600-5	120187-29-3	Muta. Cat. 3; R68 R53	Xn R: 68-53 S: (2-)22-36/37-61		
616-074-00-8	*N*-butyl-2-(4-morpholinylcarbonyl)benzamide	407-730-2	104958-67-0	Xi; R36 R43 R52-53	Xi R: 36-43-52/53 S: (2-)24-26-37-61		
616-075-00-3	D,L-(*N*,*N*-diethyl-2-hydroxy-2-phenylacetamide)	408-120-9	65197-96-8	Xn; R22 Xi; R41	Xn R: 22-41 S: (2-)26-39-46		

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616-076-00-9	tebufenozide (ISO); *N-tert*-butyl-*N'*-(4-ethylbenzoyl)-3,5-dimethylbenzohydrazide	412-850-3	112410-23-8	N; R51-53	N R: 51/53 S: 61		
616-077-00-4	reaction mass of: 2-(9-methyl-1,3,8,10-tetraoxo-2,3,9,10-tetrahydro-(1*H*,8*H*)-anthra[2,1,9-*def*: 6,5,10-*d'e'f'*]diisoquinolin-2-ylethansulfonic acid; potassium 2-(9-methyl-1,3,8,10-tetraoxo-2,3,9,10-tetrahydro-(1*H*,8*H*)-anthra[2,1,9-*def*: 6,5,10-*d'e'f'*]diisoquinolin-2-ylethansulfate	411-310-4	—	Xi; R41	Xi R: 41 S: (2-)26-39		
616-078-00-X	2-[2,4-bis(1,1-dimethyl-ethyl)phenoxy]-*N*-(2-hydroxy-5-methyl-phenyl)hexanamide	411-330-3	104541-33-5	R53	R: 53 S: 61		
616-079-00-5	1,6-hexanediyl-bis(2-(2-(1-ethylpentyl)-3-oxazolidinyl)ethyl) carbamate	411-700-4	140921-24-0	R43	Xi R: 43 S: (2-)24-37		
616-080-00-0	4-(2-((3-ethyl-4-methyl-2-oxo-pyrrolin-1-yl)carboxamido) ethyl)benzenesulfonamide)	411-850-0	119018-29-0	R52-53	R: 52/53 S: 61		
616-081-00-6	5-bromo-8-naphtholactam	413-480-5	24856-00-6	Xn; R22 R43 N; R50-53	Xn; N R: 22-43-50/53 S: (2-)22-24-37-60-61		
616-082-00-1	*N*-(5-chloro-3-((4-(diethylamino)-2-methylphenyl)imino-4-methyl-6-oxo-1,4-cyclohexadien-1-yl)benzamide	413-200-1	129604-78-0	R43	Xi R: 43 S: (2-)24-37		
616-083-00-7	[2-[(4-nitrophenyl)amino]ethyl]urea	410-700-1	27080-42-8	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
616-084-00-2	2,4-bis[*N'*-(4-methylphenyl)ureido]toluene	411-790-5	—	N; R50-53	N R: 50/53 S: 60-61		
616-085-00-8	3-(2,4-dichlorophenyl)-6-fluoro-quinazoline-2,4(1*H*,3*H*)-dione	412-190-6	168900-02-5	N; R50-53	N R: 50/53 S: 60-61		
616-086-00-3	2-acetylamino-6-chloro-4-[(4-diethylamino)2-methylphenyl-imino]-5-methyl-1-oxo-2,5-cyclohexadiene	412-250-1	102387-48-4	R53	R: 53 S: 61		

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616-087-00-9	reaction mass of: 7,9,9-trimethyl-3,14-dioxa-4,13-dioxo-5,12-diazahexadecane-1,16-diyl-prop-2-enoate; 7,7,9-trimethyl-3,14-dioxa-4,13-dioxo-5,12-diazahexadecan-1,16-diyl-prop-2-enoate	412-260-6	52658-19-2	Xi; R36 R43 N; R51-53	Xi; N R: 36-43-51/53 S: (2-)26-36/37-61		
616-088-00-4	2-aminosulfonyl-*N*,*N*-dimethylnicotinamide	413-440-7	112006-75-4	R43 R52-53	Xi R: 43-52/53 S: (2-)24-37-61		
616-089-00-X	5-(2,4-dioxo-1,2,3,4-tetrahydropyrimidine)-3-fluoro-2-hydroxymethyltetrahydrofuran	415-360-8	41107-56-6	Muta. Cat. 3; R68	Xn R: 68 S: (2-)22-36/37		
616-090-00-5	1-(1,4-benzodioxan-2-ylcarbonyl)piperazine hydrochloride	415-660-9	70918-74-0	T; R23/24/25 Xn; R48/22 N; R51-53	T; N R: 23/24/25-48/22-51/53 S: 53-45-61		
616-091-00-0	1,3,5-tris-[(2*S* and 2*R*)-2,3-epoxypropyl]-1,3,5-triazine-2,4,6-(1*H*,3*H*,5*H*)-trione	423-400-0	59653-74-6	Muta. Cat. 2; R46 T; R23 Xn; R22-48/22 Xi; R41 R43	T R: 46-22-23-41-43-48/22 S: 53-45		E
616-092-00-6	Polymeric reaction product of bicyclo[2.2.1]hepta-2,5-diene, ethene, 1,4-hexadiene, 1-propene with *N*,*N*-di-2-propenyl-formamide	404-035-6	—	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
616-093-00-1	Reaction products of: aniline-terephthalaldehyde-*o*-toluidine condensate with maleic anhydride	406-620-1	129217-90-9	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
616-094-00-7	3,3'-dicyclohexyl-1,1'-methylenebis(4,1-phenylene)diurea	406-370-3	58890-25-8	R43 R53	Xi R: 43-53 S: (2-)24-37-61		
616-095-00-2	3,3'-dioctadecyl-1,1'-methylenebis(4,1-phenylene)diurea	406-690-3	43136-14-7	R53	R: 53 S: 61		
616-096-00-8	*N*-(3-hexadecyloxy-2-hydroxyprop-1-yl)-*N*-(2-hydroxyethyl) palmitamide	408-110-4	110483-07-3	R53	R: 53 S: 61		

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616-097-00-3	*N*,*N*'-1,4-phenylenebis(2-((2-methoxy-4-nitrophenyl)azo)-3-oxobutanamide	411-840-6	83372-55-8	R53	R: 53 S: 61		
616-098-00-9	1-[4-chloro-3-((2,2,3,3,3-pentafluoropropoxy)methyl)phenyl]-5-phenyl-1*H*-1,2,4-triazole-3-carboxamide	411-750-7	119126-15-7	N; R51-53	N R: 51/53 S: 61		
616-099-00-4	2-[4-[(4-hydroxyphenyl)sulfonyl]phenoxy]-4,4-dimethyl-*N*-[5-[(methylsulfonyl)amino]-2-[4-(1,1,3,3-tetramethylbutyl)phenoxy]phenyl]-3-oxopentanamide	414-170-2	135937-20-1	R53	R: 53 S: 61		
616-100-00-8	1,3-dimethyl-1,3-bis(trimethylsilyl)urea	414-180-7	10218-17-4	Xn; R22 Xi; R38	Xn R: 22-38 S: (2-)36/37		
616-101-00-3	(*S*)-*N*-*tert*-butyl-1,2,3,4-tetrahydro-3-isoquinolinecarboxamide	414-600-9	149182-72-9	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)61		
616-102-00-9	reaction mass of: α-[3-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyl]-ω-[3-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyloxy]-poly-(oxyethylene-co-oxypropylene); 1,2-(or 1,3-)bis[α-(3-mercaptopropanoxycarbonylamino)methylphenylaminocarbonyl)-ω-oxy-poly(oxyethylene-co-oxypropylene)]-3-(or 2-)propanol; 1,2,3-tris[α-(3-mercaptopropanoxycarbonyl-amino)methylphenylaminocarbonyl)-ω-oxy-poly-(oxyethylene-co-oxypropylene)]propane]	415-870-0	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)36/37-61		
616-103-00-4	(*S*,*S*)-*trans*-4-(acetylamino)-5,6-dihydro-6-methyl-7,7-dioxo-4*H*-thieno[2,3-b]thiopyran-2-sulfonamide	415-030-3	120298-38-6	R43 N; R50-53	Xi; N R: 43-50/53 S: (2-)24-37-60-61		
616-104-00-X	benalaxyl (ISO); methyl *N*-(2,6-dimethylphenyl)-*N*-(phenylacetyl)-DL-alaninate	275-728-7	71626-11-4	N; R50-53	N R: 50/53 S: 60-61		
616-105-00-5	chlorotoluron (ISO); 3-(3-chloro-*p*-tolyl)-1,1-dimethylurea	239-592-2	15545-48-9	Carc. Cat. 3; R40 Repr. Cat. 3; R63 N; R50-53	Xn; N R: 40-63-50/53 S: (2-)26-36/37-46-60-61		

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616-106-00-0	phenmedipham (ISO); methyl 3-(3-methylcarbaniloyloxy)carbanilate	237-199-0	13684-63-4	N; R50-53	N R: 50/53 S: 60-61		
616-108-00-1	iodosulfuron-methyl-sodium; sodium ({[5-iodo-2-(methoxycarbonyl)phenyl]sulfonyl}carbamoyl)(4-methoxy-6-methyl-1,3,5-triazin-2-yl)azanide	—	144550-36-7	N; R50-53	N R: 50/53 S: 60-61		
616-109-00-7	sulfosulfuron (ISO); 1-(4,6-dimethoxypyrimidin-2-yl)-3-(2-ethylsulfonylimidazo[1,2-a]pyridin-3-yl)sulfonylurea	—	141776-32-1	N; R50-53	N R: 50/53 S: 60-61		
616-110-00-2	cyclanilide (ISO); 1-(2,4-dichloroanilinocarbonyl)cyclopropanecarboxylic acid	419-150-7	113136-77-9	Xn; R22 N; R51-53	Xn; N R: 22-51/53 S: (2-)61		
616-111-00-8	fenhexamid (ISO); *N*-(2,3-dichlor-4-hydroxyphenyl)-1-methylcyclohexancarboxamid	422-530-5	126833-17-8	N; R51-53	N R: 51/53 S: 61		
616-112-00-3	oxasulfuron (ISO); oxetan-3-yl 2-[(4,6-dimethylpyrimidin-2-yl)-carbamoylsulfamoyl]benzoate	—	144651-06-9	Xn; R48/22 N; R50-53	Xn; N R: 48/22-50/53 S: (2-)46-60-61		
616-113-00-9	desmedipham (ISO); ethyl 3-phenylcarbamoyloxyphenylcarbamate	237-198-5	13684-56-5	N; R50-53	N R: 50/53 S: 60-61	N; R50-53: C ≥ 2,5 % N; R51-53: 0,25 % ≤ C < 2,5 % R52-53: 0,025 % ≤ C < 0,25 %	
616-114-00-4	dodecanamide, *N,N'*-(9,9',10,10'-tetrahydro-9,9',10,10'-tetraoxo(1,1'-bianthracene)-4,4'-diyl)bis-	418-010-2	136897-58-0	R53	R: 53 S: 22-61		
616-115-00-X	*N*-(3-acetyl-2-hydroxyphenyl)-4-(4-phenylbutoxy)benzamide	416-150-9	136450-06-1	R53	R: 53 S: 61		
616-116-00-5	*N*-(4-dimethylaminopyridinium)-3-methoxy-4-(1-methyl-5-nitroindol-3-ylmethyl)-*N*-(*o*-tolylsulfonyl)benzamidate	416-790-9	143052-96-4	R53	R: 53 S: 61		
616-117-00-0	*N*-[2-(3-acetyl-5-nitrothiophen-2-ylazo)-5-diethylaminophenyl]acetamide	416-860-9	777891-21-1	Repr. Cat. 3; R62 R43 N; R50-53	Xn; N R: 43-62-50/53 S: (2-)22-36/37-60-61		

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616-118-00-6	*N*-(2',6'-dimethylphenyl)-2-piperidinecarboxamide hydrochloride	417-950-0	65797-42-4	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
616-119-00-1	2-(1-butyl-3,5-dioxo-2-phenyl-(1,2,4)-triazolidin-4-yl)-4,4-dimethyl-3-oxo-*N*-(2-methoxy-5-(2-(dodecyl-1-sulfonyl))propionylamino)-phenyl)-pentanamide	418-060-5	118020-93-2	R53	R: 53 S: 61		
616-120-00-7	reaction mass of: *N*-(3-dimethylamino-4-methyl-phenyl)-benzamide; *N*-(3-dimethylamino-2-methyl-phenyl)-benzamide; *N*-(3-dimethylamino-3-methyl-phenyl)-benzamide	420-600-1	—	Xn; R48/22 N; R51-53	Xn; N R: 48/22-51/53 S: (2-)36/37-61		
616-121-00-2	2,4-dihydroxy-*N*-(2-methoxyphenyl)benzamide	419-090-1	129205-19-2	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
616-123-00-3	*N*-[3-[[4-(diethylamino)-2-methylphenyl]imino]-6-oxo-1,4-cyclohexadienyl]acetamide	414-740-0	96141-86-5	N; R50-53	N R: 50/53 S: 60-61		
616-124-00-9	lithium bis(trifluoromethylsulfonyl)imide	415-300-0	90076-65-6	T; R24/25 C; R34 R52-53	T R: 24/25-34-52/53 S: (1/2-)22-26-36/37/39-45-61		
616-125-00-4	3-cyano-*N*-(1,1-dimethylethyl)androsta-3,5-diene-17-β-carboxamide	415-730-9	151338-11-3	N; R50-53	N R: 50/53 S: 60-61		
616-127-00-5	reaction mass of: *N*,*N*'-Ethane-1,2-diylbis(decanamide); 12-Hydroxy-*N*-[2-[1-oxydecyl)amino]ethyl]octadecanamide; *N*,*N*'-Ethane-1,2-diylbis(12-hydroxyoctadecanamide)	430-050-2	—	R43 N; R51-53	Xi; N R: 43-51/53 S: (2-)24-37-61		
616-128-00-0	*N*-(2-(1-allyl-4,5-dicyanoimidazol-2-ylazo)-5-(dipropylamino) phenyl)-acetamide	417-530-7	123590-00-1	R53	R: 53 S: 61		
616-129-00-6	*N*,*N*'-bis(2,2,6,6-tetramethyl-4-piperidyl)isophthalamide	419-710-0	42774-15-2	Xn; R22 Xi; R36	Xn R: 22-36 S: (2-)22-25-26		
616-130-00-1	*N*-(3-(2-(4,4-dimethyl-2,5-dioxo-imidazolin-1-yl)-4,4-dimethyl-3-oxo-pentanoylamino)-4-methoxy-phenyl)-octadecanamide	421-780-2	150919-56-5	R53	R: 53 S: 61		

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616-132-00-2	N-[4-(4-cyano-2-furfurylidene-2,5-dihydro-5-oxo-3-furyl) phenyl]butane-1-sulfonamide	423-250-6	130016-98-7	N; R50-53	N R: 50/53 S: 60-61		
616-133-00-8	*N*-cyclohexyl-*S*,*S*-dioxobenzo[*b*]tiophene-2-carboxamide	423-990-1	149118-66-1	Xn; R22 Xi; R41 N; R50-53	Xn; N R: 22-41-50/53 S: (2-)22-26-39-60-61		
616-134-00-3	3,3'-bis(dioctyloxyphosphinothioylthio)-*N*,*N*'-oxybis(methylene)dipropionamide	401-820-5	793710-14-2	R52-53	R: 52/53 S: 61		
616-135-00-9	(3*S*,4a*S*,8a*S*)-2-[(2*R*,3*S*)-3-amino-2-hydroxy-4-phenylbutyl]-*N*-tert-butyldecahydroisoquinoline-3-carboxamide	430-230-0	136522-17-3	Xn; R22 R52-53	Xn R: 22-52/53 S: (2-)22-61		
616-142-00-7	1,3-Bis(vinylsulfonylacetamido)propane	428-350-3	93629-90-4	Muta. Cat. 3; R68 Xi; R41 R43 R52-53	Xn R: 41-43-68-52/53 S: (2-)22-26-36/37/39-61		
616-143-00-2	*N*,*N*'-dihexadecyl-*N*,*N*'-bis(2-hydroxyethyl)propanediamide	422-560-9	149591-38-8	Repr. Cat. 3; R62 Xi; R36 R53	Xn R: 36-62-53 S: (2-)26-36/37-61		
617-001-00-2	di-*tert*-butyl peroxide	203-733-6	110-05-4	O; R7 F; R11	O; F R: 7-11 S: (2-)3/7-14-16-36/37/39		
617-002-00-8	α,α-dimethylbenzyl hydroperoxide; cumene hydroperoxide	201-254-7	80-15-9	O; R7 T; R23 Xn; R21/22-48/20/22 C; R34 N; R51-53	O; T; N R: 7-21/22-23-34-48/20/22-51/53 S: (1/2-)3/7-14-36/37/39-45-50-61	C; R34: C ≥ 10 % Xi; R37/38-41: 3 % ≤ C < 10 % Xi; R36/37: 1 % ≤ C < 3 %	
617-003-00-3	dilauroyl peroxide	203-326-3	105-74-8	O; R7	O R: 7 S: (2-)3/7-14-36/37/39		

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617-004-00-9	1,2,3,4-tetrahydro-1-naphthyl hydroperoxide	212-230-0	771-29-9	O; R7 Xn; R22 C; R34 N; R50-53	O; C; N R: 7-22-34-50/53 S: (1/2-)3/7-14-26-36/37/39-45-60-61	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
617-006-00-X	bis(α,α-dimethylbenzyl) peroxide	201-279-3	80-43-3	O; R7 Xi; R36/38 N; R51-53	O; Xi; N R: 7-36/38-51/53 S: (2-)3/7-14-36/37/39-61		
617-007-00-5	*tert*-butyl α,α-dimethylbenzyl peroxide	222-389-8	3457-61-2	O; R7 Xi; R38 N; R51-53	O; Xi; N R: 7-38-51/53 S: (2-)3/7-14-36/37/39-61		
617-008-00-0	dibenzoyl peroxide; benzoyl peroxide	202-327-6	94-36-0	E; R2 ⊗ Xi; R36 R43	E; Xi R: 2-36-43 S: (2-)3/7-14-36/37/39		
617-010-00-1	1-hydroperoxycyclohexyl 1-hydroxycyclohexyl peroxide; [1] 1,1'-dioxybiscyclohexan-1-ol; [2] cyclohexylidene hydroperoxide; [3] cyclohexanone, peroxide [4]	201-091-1 [1] 219-306-2 [2] 220-279-4 [3] 235-527-7 [4]	78-18-2 [1] 2407-94-5 [2] 2699-11-8 [3] 12262-58-7 [4]	E; R2 ⊗ Xn; R22 C; R34	E; C R: 2-22-34 S: (1/2-)3/7-14-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	C
617-012-00-2	8-*p*-menthyl hydroperoxide; *p*-menthane hydroperoxide	201-281-4	80-47-7	O; R7 C; R34 Xn; R20	O; C R: 7-20-34 S: (1/2-)3/7-14-36/37/39-45	C; R34: C ≥ 10 % Xi; R36/37/38: 5 % ≤ C < 10 %	
617-013-00-8	O,O-*tert*-butyl O-docosyl monoperoxyoxalate	404-300-6	116753-76-5	O; R7 N; R50-53	O; N R: 7-50/53 S: (2-)7-14-36/37/39-47-60-61		
617-014-00-3	6-(nonylamino)-6-oxo-peroxyhexanoic acid	406-680-9	104788-63-8	O; R7 Xi; R41 R43 N; R50	O; Xi; N R: 7-41-43-50 S: (2-)3/7-14-26-36/37/39-61		

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617-015-00-9	bis(4-methylbenzoyl)peroxide	407-950-9	895-85-2	E; R2 O; R7 N; R50-53	E; N R: 2-7-50/53 S: (2-)7-14-36/37/39-47-60-61		
617-016-00-4	3-hydroxy-1,1-dimethylbutyl 2-ethyl-2-methylheptaneperoxoate	413-910-1	—	O; R7 R10 Xi; R38 N; R50-53	O; Xi; N R: 7-10-38-50/53 S: (2-)7/47-14-36/37/39-60-61		
617-017-00-X	reaction mass of: 2,2'-bis(*tert*-pentylperoxy)-*p*-diisopropylbenzene; 2,2'-bis(*tert*-pentylperoxy)-*m*-diisopropylbenzene	412-140-3	32144-25-5	O; R7 ⊗ R53	O R: 7-53 S: (2-)3/7-14-36/37/39-61		
617-018-00-5	reaction mass of: 1-methyl-1-(3-(1-methylethyl)phenyl)ethyl-1-methyl-1-phenylethylperoxide, 63 % by weight; 1-methyl-1-(4-(1-methylethyl)phenyl)ethyl-1-methyl-1-phenylethylperoxide, 31 % by weight	410-840-3	71566-50-2	O; R7 N; R51-53	O; N R: 7-51/53 S: (2-)3/7-14-36/37/39-61		
617-019-00-0	6-(phthalimido)peroxyhexanoic acid	410-850-8	128275-31-0	O; R7 Xi; R41 N; R50	O; Xi; N R: 7-41-50 S: (2-)3/7-14-26-36/37/39-61		
617-020-00-6	1,3-di(prop-2,2-diyl)benzene bis(neodecanoylperoxide)	420-060-5	117663-11-3	R10 O; R7 N; R51-53	O; N R: 7-10-51/53 S: (2-)7-14-36/37/39-47-61		
647-001-00-8	glucosidase, β-	232-589-7	9001-22-3	R42	Xn R: 42 S: (2-)22-24-36/37		
647-002-00-3	cellulase	232-734-4	9012-54-8	R42	Xn R: 42 S: (2-)22-24-36/37		
647-003-00-9	cellobiohydrolase, exo-	253-465-9	37329-65-0	R42	Xn R: 42 S: (2-)22-24-36/37		

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647-004-00-4	cellulases with the exception of those specified elsewhere in this Annex	—	—	R42	Xn R: 42 S: (2-)22-24-36/37		A
647-005-00-X	bromelain, juice	232-572-4	9001-00-7	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-006-00-5	ficin	232-599-1	9001-33-6	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-007-00-0	papain	232-627-2	9001-73-4	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-008-00-6	pepsin A	232-629-3	9001-75-6	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-009-00-1	rennin	232-645-0	9001-98-3	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-010-00-7	trypsin	232-650-8	9002-07-7	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-011-00-2	chymotrypsin	232-671-2	9004-07-3	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-012-00-8	subtilisin	232-752-2	9014-01-1	Xi; R37/38-41 R42	Xn R: 37/38-41-42 S: (2-)22-24-26-36/37/39		
647-013-00-3	proteinase, microbial neutral	232-966-6	9068-59-1	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		

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647-014-00-9	proteases with the exception of those specified elsewhere in this Annex	—	—	Xi; R36/37/38 R42	Xn R: 36/37/38-42 S: (2-)22-24-26-36/37		
647-015-00-4	amylase, α-	232-565-6	9000-90-2	R42	Xn R: 42 S: (2-)22-24-36/37		
647-016-00-X	amylases with the exception of those specified elsewhere in this Annex	—	—	R42	Xn R: 42 S: (2-)22-24-36/37		
648-001-00-0	Distillates (coal tar), benzole fraction; Light Oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists of hydrocarbons having carbon numbers primarily in the range of C_4 to C_{10} and distilling in the approximate range of 80 °C to 160 °C (175 °F to 320 °F).]	283-482-7	84650-02-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-002-00-6	Tar oils, brown-coal; Light Oil; [The distillate from lignite tar boiling in the range of approximately 80 °C to 250 °C (176 °F to 482 °F). Composed primarily of aliphatic and aromatic hydrocarbons and monobasic phenols.]	302-674-4	94114-40-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-003-00-1	Benzol forerunnings (coal); Light Oil Redistillate, low boiling; [The distillate from coke oven light oil having an approximate distillation range below 100 °C (212 °F). Composed primarily of C_4 to C_6 aliphatic hydrocarbons.]	266-023-5	65996-88-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-004-00-7	Distillates (coal tar), benzole fraction, BTX-rich; Light Oil Redistillate, low boiling; [A residue from the distillation of crude benzole to remove benzole fronts. Composed primarily of benzene, toluene and xylenes boiling in the range of approximately 75 °C to 200 °C (167 °F to 392 °F).]	309-984-9	101896-26-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-005-00-2	Aromatic hydrocarbons, C_{6-10}, C_8-rich; Light Oil Redistillate, low boiling	292-697-5	90989-41-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

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648-006-00-8	Solvent naphtha (coal), light; Light Oil Redistillate, low boiling	287-498-5	85536-17-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-007-00-3	Solvent naphtha (coal), xylene-styrene cut; Light Oil Redistillate, intermediate boiling	287-502-5	85536-20-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-008-00-9	Solvent naphtha (coal), coumarone-styrene contg.; Light Oil Redistillate, intermediate boiling	287-500-4	85536-19-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-009-00-4	Naphtha (coal), distn. residues; Light Oil Redistillate, high boiling; [The residue remaining from the distillation of recovered naphtha. Composed primarily of naphthalene and condensation products of indene and styrene.]	292-636-2	90641-12-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-010-00-X	Aromatic hydrocarbons, C_8; Light Oil Redistillate, high boiling	292-694-9	90989-38-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-012-00-0	Aromatic hydrocarbons, C_{8-9}, hydrocarbon resin polymn. by-product; Light Oil Redistillate, high boiling; [A complex combination of hydrocarbons obtained from the evaporation of solvent under vacuum from polymerized hydrocarbon resin. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_9 and boiling in the range of approximately 120 °C to 215 °C (248 °F to 419 °F).]	295-281-1	91995-20-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-013-00-6	Aromatic hydrocarbons, C_{9-12}, benzene distn.; Light Oil Redistillate, high boiling	295-551-9	92062-36-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-014-00-1	Extract residues (coal), benzole fraction alk., acid ext.; Light Oil Extract Residues, low boiling; [The redistillate from the distillate, freed of tar acids and tar bases, from bituminous coal high temperature tar boiling in the approximate range of 90 °C to 160 °C (194 °F to 320 °F). It consists predominantly of benzene, toluene and xylenes.]	295-323-9	91995-61-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-015-00-7	Extract residues (coal tar), benzole fraction alk., acid ext.; Light Oil Extract Residues, low boiling; [A complex combination of hydrocarbons obtained by the redistillation of the distillate of high temperature coal tar (tar acid and tar base free). It consists predominantly of unsubstituted and substituted mononuclear aromatic hydrocarbons boiling in the range of 85 °C-195 °C (185 °F-383 °F).]	309-868-8	101316-63-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-016-00-2	Extract residues (coal), benzole fraction acid; Light Oil Extract Residues, low boiling; [An acid sludge by-product of the sulphuric acid refining of crude high temperature coal. Composed primarily of sulfuric acid and organic compounds.]	298-725-2	93821-38-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-017-00-8	Extract residues (coal), light oil alk., distn. overheads; Light Oil Extract Residues, low boiling; [The first fraction from the distillation of aromatic hydrocarbons, coumarone, naphthalene and indene rich prefactionator bottoms or washed carbolic oil boiling substantially below 145 °C (293 °F). Composed primarily of C_7 and C_8 aliphatic and aromatic hydrocarbons.]	292-625-2	90641-02-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-018-00-3	Extract residues (coal), light oil alk., acid ext., indene fraction; Light Oil Extract Residues, intermediate boiling	309-867-2	101316-62-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-019-00-9	Extract residues (coal), light oil alk., indene naphtha fraction; Light Oil Extract Residues, high boiling; [The distillate from aromatic hydrocarbons, coumarone, naphthalene and indene rich prefractionator bottoms or washed carbolic oils, having an approximate boiling range of 155 °C to 180 °C (311 °F to 356 °F). Composed primarily of indene, indan and trimethylbenzenes.]	292-626-8	90641-03-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-020-00-4	Solvent naphtha (coal); Light Oil Extract Residues, high boiling; [The distillate from either high temperature coal tar, coke oven light oil, or coal tar oil alkaline extract residue having an approximate distillation range of 130 °C to 210 °C (266 °F to 410 °F) Composed primarily of indene and other polycyclic ring systems containing a single aromatic ring. May contain phenolic compounds and aromatic nitrogen bases.]	266-013-0	65996-79-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

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648-021-00-X	Distillates (coal tar), light oils, neutral fraction; Light Oil Extract Residues, high boiling; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of alkyl-substituted one ring aromatic hydrocarbons boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F). May also include unsaturated hydrocarbons such as indene and coumarone.]	309-971-8	101794-90-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-022-00-5	Distillates (coal tar), light oils, acid exts.; Light Oil Extract Residues, high boiling; [This oil is a complex mixture of aromatic hydrocarbons, primarily indene, naphthalene, coumarone, phenol, and *o*-, *m*- and *p*-cresol and boiling in the range of 140 °C to 215 °C (284 °F to 419 °F).]	292-609-5	90640-87-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-023-00-0	Distillates (coal tar), light oils; Carbolic Oil; [A complex combination of hydrocarbons obtained by distillation of coal tar. It consists of aromatic and other hydrocarbons, phenolic compounds and aromatic nitrogen compounds and distills at the approximate range of 150 °C to 210 °C (302 °F to 410 °F).]	283-483-2	84650-03-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-024-00-6	Tar oils, coal; Carbolic Oil; [The distillate from high temperature coal tar having an approximate distillation range of 130 °C to 250 °C (266 °F to 410 °F). Composed primarily of naphthalene, alkylnaphthalenes, phenolic compounds, and aromatic nitrogen bases.]	266-016-7	65996-82-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-026-00-7	Extract residues (coal), light oil alk., acid ext.; Carbolic Oil Extract Residue; [The oil resulting from the acid washing of alkali-washed carbolic oil to remove the minor amounts of basic compounds (tar bases). Composed primarily of indene, indan and alkylbenzenes.]	292-624-7	90641-01-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-027-00-2	Extract residues (coal), tar oil alk.; Carbolic Oil Extract Residue; [The residue obtained from coal tar oil by an alkaline wash such as aqueous sodium hydroxide after the removal of crude coal tar acids. Composed primarily of naphthalenes and aromatic nitrogen bases.]	266-021-4	65996-87-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-028-00-8	Extract oils (coal), light oil; Acid Extract; [The aqueous extract produced by an acidic wash of alkali-washed carbolic oil. Composed primarily of acid salts of various aromatic nitrogen bases including pyridine, quinoline and their alkyl derivatives.]	292-622-6	90640-99-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-029-00-3	Pyridine, alkyl derivs.; Crude Tar Bases; [The complex combination of polyalkylated pyridines derived from coal tar distillation or as high-boiling distillates approximately above 150 °C (302 °F) from the reaction of ammonia with acetaldehyde, formaldehyde or paraformaldehyde.]	269-929-9	68391-11-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-030-00-9	Tar bases, coal, picoline fraction; Distillate Bases; [Pyridine bases boiling in the range of approximately 125 °C to 160 °C (257 °F 320 °F) obtained by distillation of neutralized acid extract of the base-containing tar fraction obtained by the distillation of bituminous coal tars. Composed chiefly of lutidines and picolines.]	295-548-2	92062-33-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-031-00-4	Tar bases, coal, lutidine fraction; Distillate Bases	293-766-2	91082-52-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-032-00-X	Extract oils (coal), tar base, collidine fraction; Distillate Bases; [The extract produced by the acidic extraction of bases from crude coal tar aromatic oils, neutralization, and distillation of the bases. Composed primarily of collidines, aniline, toluidines, lutidines, xylidines.]	273-077-3	68937-63-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-033-00-5	Tar bases, coal, collidine fraction; Distillate Bases; [The distillation fraction boiling in the range of approximately 181 °C to 186 °C (356 °F to 367 °F) from the crude bases obtained from the neutralized, acid-extracted base-containing tar fractions obtained by the distillation of bituminous coal tar. It contains chiefly aniline and collidines.]	295-543-5	92062-28-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-034-00-0	Tar bases, coal, aniline fraction; Distillate Bases; [The distillation fraction boiling in the range of approximately 180 °C to 200 °C (356 °F to 392 °F) from the crude bases obtained by dephenolating and debasing the carbolated oil from the distillation of coal tar. It contains chiefly aniline, collidines, lutidines and toluidines.]	295-541-4	92062-27-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-035-00-6	Tar bases, coal, toluidine fraction; Distillate Bases	293-767-8	91082-53-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-036-00-1	Distillates (petroleum), alkene-alkyne manuf. pyrolysis oil, mixed with high-temp. coal tar, indene fraction; Redistillates; [A complex combination of hydrocarbons obtained as a redistillate from the fractional distillation of bituminous coal high temperature tar and residual oils that are obtained by the pyrolytic production of alkenes and alkynes from petroleum products or natural gas. It consists predominantly of indene and boils in a range of approximately 160 °C to 190 °C (320 °F to 374 °F).]	295-292-1	91995-31-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-037-00-7	Distillates (coal), coal tar-residual pyrolysis oils, naphthalene oils; Redistillates; [The redistillate obtained from the fractional distillation of bituminous coal high temperature tar and pyrolysis residual oils and boiling in the range of approximately 190 °C to 270 °C (374 °F to 518 °F). Composed primarily of substituted dinuclear aromatics.]	295-295-8	91995-35-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-038-00-2	Extract oils (coal), coal tar-residual pyrolysis oils, naphthalene oil, redistillate; Redistillates; [The redistillate from the fractional distillation of dephenolated and debased methylnaphthalene oil obtained from bituminous coal high temperature tar and pyrolysis residual oils boiling in the approximate range of 220 °C to 230 °C (428 °F to 446 °F). It consists predominantly of unsubstituted and substituted dinuclear aromatic hydrocarbons.]	295-329-1	91995-66-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-039-00-8	Extract oils (coal), coal tar-residual pyrolysis oils, naphthalene oils; Redistillates; [A neutral oil obtained by debasing and dephenolating the oil obtained from the distillation of high temperature tar and pyrolysis residuel oils which has a boiling range of 225 °C to 255 °C (437 °F to 491 °F). Composed primarily of substituted dinuclear aromatic hydrocarbons.]	310-170-0	122070-79-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-040-00-3	Extract oils (coal), coal tar residual pyrolysis oils, naphthalene oil, distn. residues; Redistillates; [Residue from the distillation of dephenolated and debased methylnaphthalene oil (from bituminous coal tar and pyrolysis residual oils) with a boiling range of 240 °C to 260 °C (464 °F to 500 °F). Composed primarily of substituted dinuclear aromatic and heterocyclic hydrocarbons.]	310-171-6	122070-80-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-041-00-9	Absorption oils, bicyclo arom. and heterocyclic hydrocarbon fraction; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained as a redistillate from the distillation of wash oil. It consists predominantly of 2-ringed aromatic and heterocyclic hydrocarbons boiling in the range of approximately 260 °C to 290 °C (500 °F to 554 °F).]	309-851-5	101316-45-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-042-00-4	Distillates (coal tar), upper, fluorene-rich; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained by the crystallization of tar oil. It consists af aromatic and polycyclic hydrocarbons primarily fluorene and some acenaphthene.]	284-900-0	84989-11-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-043-00-X	Creosote oil, acenaphthene fraction, acenaphthene-free; Wash Oil Redistillate; [The oil remaining after removal by a crystallization process of acenaphthene from acenaphthene oil from coal tar. Composed primarily of naphthalene and alkylnaphthalenes.]	292-606-9	90640-85-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-044-00-5	Distillates (coal tar), heavy oils; Heavy Anthracene Oil; [Distillate from the fractional distillation of coal tar of bituminous coal, with boiling range of 240 °C to 400 °C (464 °F to 752 °F). Composed primarily of tri- and polynuclear hydrocarbons and heterocyclic compounds.]	292-607-4	90640-86-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-045-00-0	Distillates (coal tar), upper; Heavy Anthracene Oil; [The distillate from coal tar having an approximate distillation range of 220 °C to 450 °C (428 °F to 842 °F). Composed primarily of three to four membered condensed ring aromatic hydrocarbons and other hydrocarbons.]	266-026-1	65996-91-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-046-00-6	Anthracene oil, acid ext.; Anthracene Oil Extract Residue; [A complex combination of hydrocarbons from the base-freed fraction obtained from the distillation of coal tar and boiling in the range of approximately 325 °C to 365 °C (617 °F to 689 °F). It contains predominantly anthracene and phenanthrene and their alkyl derivatives.]	295-274-3	91995-14-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-047-00-1	Distillates (coal tar); Heavy Anthracene Oil; [The distillate from coal tar having an approximate distillation range of 100 °C to 450 °C (212 °F to 842 °F). Composed primarily of two to four membered condensed ring aromatic hydrocarbons, phenolic compounds, and aromatic nitrogen bases.]	266-027-7	65996-92-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-048-00-7	Distillates (coal tar), pitch, heavy oils; Heavy Anthracene Oil; [The distillate from the distillation of the pitch obtained from bituminous high temperature tar. Composed primarily of tri- and polynuclear aromatic hydrocarbons and boiling in the range of approximately 300 °C to 470 °C (572 °F to 878 °F). The product may also contain heteroatoms.]	295-312-9	91995-51-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-049-00-2	Distillates (coal tar), pitch; Heavy Anthracene Oil; [The oil obtained from condensation of the vapors from the heat treatment of pitch. Composed primarily of two- to four-ring aromatic compounds boiling in the range of 200 °C to greater than 400 °C (392 °F to greater than 752 °F).]	309-855-7	101316-49-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-050-00-8	Distillates (coal tar), heavy oils, pyrene fraction; Heavy Anthracene Oil Redistillate; [The redistillate obtained from the fractional distillation of pitch distillate boiling in the range of approximately 350 °C to 400 °C (662 °F to 752 °F). Consists predominantly of tri- and polynuclear aromatics and heterocyclic hydrocarbons.]	295-304-5	91995-42-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-051-00-3	Distillates (coal tar), pitch, pyrene fraction; Heavy Anthracene Oil Redistillate; [The redistillate obtained from the fractional distillation of pitch distillate and boiling in the range of approximately 380 °C to 410 °C (7160 to 770 °F). Composed primarily of tri- and polynuclear aromatic hydrocarbons and heterocyclic compounds.]	295-313-4	91995-52-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-052-00-9	Paraffin waxes (coal), brown-coal high-temp. tar, carbon-treated; Coal Tar Extract; [A complet combination of hydrocarbons obtained by the treatment of lignite carbonization tar with activated carbon for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-296-6	97926-76-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-053-00-4	Paraffin waxes (coal), brown-coal high-temp tar, clay-treated; Coal Tar Extract; [A complex combination of hydrocarbons obtained by the treatment of lignite carbonization tar with bentonite for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-297-1	97926-77-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-054-00-X	Pitch; Pitch	263-072-4	61789-60-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-055-00-5	Pitch, coal tar, high-temp.; Pitch; [The residue from the distillation of high temperature coal tar. A black solid with an approximate softening point from 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of a complex mixture of three or more membered condensed ring aromatic hydrocarbons.]	266-028-2	65996-93-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-056-00-0	Pitch, coal tar, high-temp., heat-treated; Pitch; [The heat treated residue from the distillation of high temperature coal tar. A black solid with an approximate softening point from 80 °C to 180 °C (176 °F to 356 °F). Composed primarily of a complex mixture of three or more membered condensed ring aromatic hydrocarbons.]	310-162-7	121575-60-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-057-00-6	Pitch, coal tar, high-temp., secondary; Pitch Redistillate; [The residue obtained during the distillation of high boiling fractions from bituminous coal high temperature tar and/or pitch coke oil, with a softening point of 140 °C to 170 °C (284 °F to 392 °F) according to DIN 52025. Composed primarily of tri- and polynuclear aromatic compounds which also contain heteroatoms.]	302-650-3	94114-13-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-058-00-1	Residues (coal tar), pitch distn.; Pitch Redistillate; [Residue from the fractional distillation of pitch distillate boiling in the range of approximately 400 °C to 470 °C (752 °F to 846 °F). Composed primarily of polynuclear aromatic hydrocarbons, and heterocyclic compounds.]	295-507-9	92061-94-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-059-00-7	Tar, coal, high-temp., distn. and storage residues; Coal Tar Solids Residue; [Coke- and ash-containing solid residues that separate on distillation and thermal treatment of bituminous coal high temperature tar in distillation installations and storage vessels. Consists predominantly of carbon and contains a small quantity of hetero compounds as well as ash components.]	295-535-1	92062-20-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-060-00-2	Tar, coal, storage residues; Coal Tar Solids Residue; [The deposit removed from crude coal tar storages. Composed primarily of coal tar and carbonaceous particulate matter.]	293-764-1	91082-50-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-061-00-8	Tar, coal, high-temp., residues; Coal Tar Solids Residue; [Solids formed during the coking of bituminous coal to produce crude bituminous coal high temperature tar. Composed primarily of coke and coal particles, highly aromatized compounds and mineral substances.]	309-726-5	100684-51-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-062-00-3	Tar, coal, high-temp., high-solids; Coal Tar Solids Residue; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in the high temperature (greater than 700 °C (1292 °F)) destructive distillation of coal. Composed primarily of a complex mixture of condensed ring aromatic hydrocarbons with a high solid content of coal-type materials.]	273-615-7	68990-61-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-063-00-9	Waste solids, coal-tar pitch coking; Coal Tar Solids Residue; [The combination of wastes formed by the coking of bituminous coal tar pitch. It consists predominantly of carbon.]	295-549-8	92062-34-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-064-00-4	Extract residues (coal), brown; Coal Tar Extract; [The residue from extraction of dried coal.]	294-285-0	91697-23-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-065-00-X	Paraffin waxes (coal), brown-coal-high-temp. tar; Coal Tar Extract; [A complex combination of hydrocarbons obtained from lignite carbonization tar by solvent crystallisation (solvent deoiling), by sweating or an adducting process. It consists predominantly of straight and branched chain saturated hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-454-1	92045-71-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-066-00-5	Paraffin waxes (coal), brown-coal-high-temp. tar, hydrotreated; Coal Tar Extract; [A complex combination of hydrocarbons obtained from lignite carbonization tar by solvent crystallisation (solvent deoiling), by sweating or an adducting process treated with hydrogen in the presence of a catalyst. It consists predominantly of straight and branched chain saturated hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-455-7	92045-72-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-067-00-0	Paraffin waxes (coal), brown-coal high-temp tar, silicic acid-treated; Coal Tar Extract; [A complex combination of hydrocarbons obtained by the treatment of lignite carbonization tar with silicic acid for removal of trace constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-298-7	97926-78-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-068-00-6	Tar, coal, low-temp., distn. residues; Tar Oil, intermediate boiling; [Residues from fractional distillation of low temperature coal tar to remove oils that boil in a range up to approximately 300 °C (572 °F). Composed primarily of aromatic compounds.]	309-887-1	101316-85-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-069-00-1	Pitch, coal tar, low-temp; Pitch Residue; [A complex black solid or semi-solid obtained from the distillation of a low temperature coal tar. It has a softening point within the approximate range of 40 °C to 180 °C (104 °F to 356 °F). Composed primarily of a complex mixture of hydrocarbons.]	292-651-4	90669-57-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-070-00-7	Pitch, coal tar, low-temp., oxidized; Pitch Residue, oxidised; [The product obtained by air-blowing, at elevated temperature, low-temperature coal tar pitch. It has a softening-point within the approximate range of 70 °C to 180 °C (158 °F to 356 °F). Composed primarily of a complex mixture of hydrocarbons.]	292-654-0	90669-59-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-071-00-2	Pitch, coal tar, low-temp., heat-treated; Pitch Residue, oxidised; Pitch Residue, heat-treated; [A complex black solid obtained by the heat treatment of low temperature coal tar pitch. It has a softening point within the approximate range of 50 °C to 140 °C (122 °F to 284 °F). Composed primarily of a complex mixture of aromatic compounds.]	292-653-5	90669-58-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-072-00-8	Distillates (coal-petroleum), condensed-ring arom; Distillates; [The distillate from a mixture of coal and tar and aromatic petroleum streams having an approximate distillation range of 220 °C to 450 °C (428 °F to 842 °F). Composed primarily of 3- to 4-membered condensed ring aromatic hydrocarbons.]	269-159-3	68188-48-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-073-00-3	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polyethylene-polypropylene pyrolysis-derived; Pyrolysis Products; [A complex combination hydrocarbons obtained from mixed coal tar pitch-polyethylene-polypropylene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-956-6	101794-74-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-074-00-9	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polyethylene pyrolysis-derived; Pyrolysis Products; [A complex combination of hydrocarbons obtained from mixed coal tar pitch-polyethylene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-957-1	101794-75-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-075-00-4	Aromatic hydrocarbons, C_{20-28}, polycyclic, mixed coal-tar pitch-polystyrene pyrolysis-derived; Pyrolysis Products; [A complex combination of hydrocarbons obtained from mixed coal tar pitch-polystyrene pyrolysis. Composed primarily of polycyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{28} and having a softening point of 100 °C to 220 °C (212 °F to 428 °F) according to DIN 52025.]	309-958-7	101794-76-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-076-00-X	Pitch, coal tar-petroleum; Pitch Residues; [The residue from the distillation of a mixture of coal tar and aromatic petroleum streams. A solid with a softening point from 40 °C to 180 °C (140 °F to 356 °F). Composed primarily of a complex combination of three or more membered condensed ring aromatic hydrocarbons.]	269-109-0	68187-57-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-077-00-5	Phenanthrene, distn. residues; Heavy Anthracene Oil Redistillate; [Residue from the distillation of crude phenanthrene boiling in the approximate range of 340 °C to 420 °C (644 °F to 788 °F). It consists predominantly of phenanthrene, anthracene and carbazole.]	310-169-5	122070-78-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-078-00-0	Distillates (coal tar), upper, fluorene-free; Wash Oil Redistillate; [A complex combination of hydrocarbons obtained by the crystallization of tar oil. It consists of aromatic polycyclic hydrocarbons, primarily diphenyl, dibenzofuran and acenaphthene.]	284-899-7	84989-10-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-079-00-6	Anthracene oil; Anthracene oil; [A complex combination of polycyclic aromatic hydrocarbons obtained from coal tar having an approximate distillation range of 300 °C ot 400 °C (572 °F to 752 °F). Composed primarily of phenanthrene, anthracene and carbazole.]	292-602-7	90640-80-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-080-00-1	Residues (coal tar), creosote oil distn.; Wash Oil Redistillate; [The residue from the fractional distillation of wash oil boiling in the approximate range of 270 °C to 330 °C (518 °F to 626 °F). It consists predominantly of dinuclear aromatic and heterocyclic hydrocarbons.]	295-506-3	92061-93-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-081-00-7	Tar, coal; Coal tar; [The by-product from the destructive distillation of coal. Almost black semisolid. A complex combination of aromatic hydro-carbons, phenolic compounds, nitrogen bases and thiophene.]	232-361-7	8007-45-2	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
648-082-00-2	Tar, coal, high-temp.; Coal tar; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in the high temperature (greater than 700 °C (1292 °F)) destructive distillation of coal. A black viscous liquid denser than water. Composed primarily of a complex mixture of condensed ring aromatic hydrocarbons. May contain minor amounts of phenolic compounds and aromatic nitrogen bases.]	266-024-0	65996-89-6	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
648-083-00-8	Tar, coal, low-temp.; Coal oil; [The condensation product obtained by cooling, to approximately ambient temperature, the gas evolved in low temperature (less than 700 °C (1292 °F)) destructive distillation of coal. A black viscous liquid denser than water. Composed primarily of condensed ring aromatic hydrocarbons, phenolic compounds, aromatic nitrogen bases, and their alkyl derivatives.]	266-025-6	65996-90-9	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
648-084-00-3	Distillates (coal), coke-oven light oil, naphthalene cut; Naphthalene Oil; [The complex combination of hydrocarbons obtained from prefractionation (continuous distillation) of coke oven light oil. It consists predominantly of naphthalene, coumarone and indene and boils above 148 °C (298 °F).]	285-076-5	85029-51-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-085-00-9	Distillates (coal tar), naphthalene oils; Naphthalene Oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists primarily of aromatic and other hydrocarbons, phenolic compounds and aromatic nitrogen compounds and distills in the approximate range of 200 °C to 250 °C (392 °F to 482 °F).]	283-484-8	84650-04-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-086-00-4	Distillates (coal tar), naphthalene oils, naphthalene-low; Naphthalene Oil Redistillate; [A complex combination of hydrocarbons obtained by crystallization of naphthalene oil. Composed primarily of naphthalene, alkyl naphthalenes and phenolic compounds.]	284-898-1	84989-09-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-087-00-X	Distillates (coal tar), naphthalene oil crystn. mother liquor; Naphthalene Oil Redistillate; [A complex combination of organic compounds obtained as a filtrate from the crystallization of the naphthalene fraction from coal tar and boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). Contains chiefly naphthalene, thionaphthene and alkylnaphthalenes.]	295-310-8	91995-49-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-088-00-5	Extract residues (coal), naphthalene oil, alk.; Naphthalene Oil Extract Residue; [A complex combination of hydrocarbons obtained from the alkali washing of naphthalene oil to remove phenolic compounds (tar acids). It is composed of naphthalene and alkyl naphthalenes.]	310-166-9	121620-47-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-089-00-0	Extract residues (coal), naphthalene oil, alk., naphthalene-low; Naphthalene Oil Extract Residue; [A complex combination of hydrocarbons remaining after the removal of naphthalene from alkali-washed naphthalene oil by a crystallization process. It is composed primarily of naphthalene and alkyl naphthalenes.]	310-167-4	121620-48-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-090-00-6	Distillates (coal tar), naphthalene oils, naphthalene-free, alk. exts.; Naphthalene Oil Extract Residue; [The oil remaining after the removal of phenolic compounds (tar acids) from drained naphthalene oil by an alkali wash. Composed primarily of naphthalene and alkyl naphthalenes.]	292-612-1	90640-90-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-091-00-1	Extract residues (coal), naphthalene oil alk., distn. overheads; Naphthalene Oil Extract Residue; [The distillation from alkali-washed naphthalene oil having an approximate distillation range of 180 °C to 220 °C (356 °F to 428 °F). Composed primarily of naphthalene, alkylbenzenes, indene and indan.]	292-627-3	90641-04-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-092-00-7	Distillates (coal tar), naphthalene oils, methylnaphthalene fraction; Methylnaphthalene Oil; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of substituted two ring aromatic hydrocarbons and aromatic nitrogen bases boiling in the range of approximately 225 °C to 255 °C (437 °F to 491 °F).]	309-985-4	101896-27-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-093-00-2	Distillates (coal tar), naphthalene oils, indole-methylnaphthalene fraction; Methylnaphthalene Oil; [A distillate from the fractional distillation of high temperature coal tar. Composed primarily of indole and methylnaphthalene boiling in the range of approximately 235 °C to 255 °C (455 °F to 491 °F).]	309-972-3	101794-91-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-094-00-8	Distillates (coal tar), naphthalene oils, acid exts.; Methylnaphthalene Oil Extract Residue; [A complex combination of hydrocarbons obtained by debasing the methylnaphthalene fraction obtained by the distillation of coal tar and boiling in the range of approximately 230 °C to 255 °C (446 °F to 491 °F). Contains chiefly 1(2)-methylnaphthalene, naphthalene, dimethylnaphthalene and biphenyl.]	295-309-2	91995-48-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-095-00-3	Extract residues (coal), naphthalene oil alk., distn. residues; Methylnaphthalene Oil Extract Residue; [The residue from the distillation of alkali-washed naphthalene oil having an approximate distillation range of 220 °C to 300 °C (428 °F to 572 °F). Composed primarily of naphthalene, alkylnaphthalenes and aromatic nitrogen bases.]	292-628-9	90641-05-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-096-00-9	Extract oils (coal), acidic, tar-base free; Methylnaphthalene Oil Extract Residue; [The extract oil boiling in the range of approximately 220 °C to 265 °C (428 °F to 509 °F) from coal tar alkaline extract residue produced by an acidic wash such as aqueous sulfuric acid after distillation to remove tar bases. Composed primarily of alkylnaphthalenes.]	284-901-6	84989-12-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-097-00-4	Distillates (coal tar), benzole fraction, distn. residues; Wash Oil; [A complex combination of hydrocarbons obtained from the distillation of crude benzole (high temperature coal tar). It may be a liquid with the approximate distillation range of 150 °C to 300 °C (302 °F to 572 °F) or a semi-solid or solid with a melting point up to 70 °C (158 °F). It is composed primarily of naphthalene and alkyl naphthalenes.]	310-165-3	121620-46-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-098-00-X	Creosote oil, acenaphthene fraction; Wash Oil; [A complex combination of hydrocarbons produced by the distillation of coal tar and boiling in the range of approximately 240 °C to 280 °C (464 °F to 536 °F). Composed primarily of acenaphthene, naphthalene and alkyl naphthalene.]	292-605-3	90640-84-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-099-00-5	Creosote oil; [A complex combination of hydrocarbons obtained by the distillation of coal tar. It consists primarily of aromatic hydrocarbons and may contain appreciable quantities of tar acids and tar bases. It distills at the approximate range of 200 °C to 325 °C (392 °F to 617 °F).]	263-047-8	61789-28-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-100-00-9	Creosote oil, high-boiling distillate; Wash Oil; [The high-boiling distillation fraction obtained from the high temperature carbonization of bituminous coal which is further refined to remove excess crystalline salts. It consists primarily of creosote oil with some of the normal polynuclear aromatic salts, which are components of coal tar distillates, removed. It is crystal free at approximately 5 °C (41 °F).]	274-565-9	70321-79-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-101-00-4	Creosote; [The distillate of coal tar produced by the high temperature carbonization of bituminous coal. It consists primarily of aromatic hydrocarbons, tar acids and tar bases.]	232-287-5	8001-58-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-102-00-X	Extract residues (coal), creosote oil acid; Wash Oil Extract Residue; [A complex combination of hydrocarbons from the base-freed fraction from the distillation of coal tar, boiling in the range of approximately 250 °C to 280 °C (482 °F to 536 °F). It consists predominantly of biphenyl and isomeric diphenyl-naphthalenes.]	310-189-4	122384-77-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-103-00-5	Anthracene oil, anthracene paste; Anthracene Oil Fraction; [The anthracene-rich solid obtained by the crystallization and centrifuging of anthracene oil. It is composed primarily of anthracene, carbazole and phenanthrene.]	292-603-2	90640-81-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-104-00-0	Anthracene oil, anthracene-low; Anthracene Oil Fraction; [The oil remaining after the removal, by a crystallization process, of an anthracene-rich solid (anthracene paste) from anthracene oil. It is composed primarily of two, three and four membered aromatic compounds.]	292-604-8	90640-82-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-105-00-6	Residues (coal tar), anthracene oil distn.; Anthracene Oil Fraction; [The residue from the fraction distillation of crude anthracene boiling in the approximate range of 340 °C to 400 °C (644 °F to 752 °F). It consists predominantly of tri- and polynuclear aromatic and heterocyclic hydrocarbons.]	295-505-8	92061-92-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-106-00-1	Anthracene oil, anthracene paste, anthracene fraction; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by the crystallization of anthracene oil from bituminous high temperature tar and boiling in the range of 330 °C to 350 °C (626 °F to 662 °F). It contains chiefly anthracene, carbazole and phenanthrene.]	295-275-9	91995-15-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-107-00-7	Anthracene oil, anthracene paste, carbazole fraction; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by crystallization of anthrancene oil from bituminous coal high temperature tar and boiling in the approximate range of 350 °C to 360 °C (662 °F to 680 °F). It contains chiefly anthacene, carbazole and phenanthrene.]	295-276-4	91995-16-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-108-00-2	Anthracene oil, anthracene paste, distn. lights; Anthracene Oil Fraction; [A complex combination of hydrocarbons from the distillation of anthracene obtained by crystallization of anthracene oil from bituminous light temperature tar and boiling in the range of approximately 290 °C to 340 °C (554 °F to 644 °F). It contains chiefly trinuclear aromatics and their dihydro derivatives.]	295-278-5	91995-17-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-109-00-8	Tar oils, coal, low-temp.; Tar Oil, high boiling; [A distillate from low-temperature coal tar. Composed primarily of hydrocarbons, phenolic compounds and aromatic nitrogen bases boiling in the range of approximately 160 °C to 340 °C (320 °F to 644 °F).]	309-889-2	101316-87-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-110-00-3	Extract residues (coal), low temp. coal atar alk.; [The residue from low temperature coal tar oils after an alkaline wash, such as aqueous sodium hydroxide, to remove crude coal tar acids. Composed primarily of hydrocarbons and aromatic nitrogen bases.]	310-191-5	122384-78-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-111-00-9	Phenols, ammonia liquor ext.; Alkaline Extract; [The combination of phenols extracted, using isobutyl acetate, from the ammonia liquor condensed from the gas evolved in low-temperature (less than 700 °C (1292 °F)) destructive distillation of coal. It consists predominantly of a mixture of monohydric and dihydric phenols.]	284-881-9	84988-93-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-112-00-4	Distillates (coal tar), light oils, alk. exts.; Alkaline Extract; [The aqueous extract from carbolic oil produced by an alkaline wash such as aqueous sodium hydroxide. Composed primarily of the alkali salts of various phenolic compounds.]	292-610-0	90640-88-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-113-00-X	Extracts, coal tar oil alk.; Alkaline Extract; [The extract from coal tar oil produced by an alkaline wash such as aqueous sodium hydroxide. Composed primarily of the alkali salts of various phenolic compounds.]	266-017-2	65996-83-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-114-00-5	Distillates (coal tar), naphthalene oils, alk. exts.; Alkaline Extract; [The aqueous extract from naphthalene oil produced by an alkaline wash such as aqueous sodium hydroxid. Composed primarily of the alkali salts of various phenolic compounds.]	292-611-6	90640-89-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-115-00-0	Extract residues (coal), tar oil alk., carbonated, limed; Crude Phenols; [The product obtained by treatment of coal tar oil alkaline extract with CO_2 and CaO. Composed primarily of $CaCO_3$, $Ca(OH)_2$, Na_2CO_3 and other organic and inorganic impurities.]	292-629-4	90641-06-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-116-00-6	Tar acids, coal, crude; Crude Phenols; [The reaction product obtained by neutralizing coal tar oil alkaline extract with an acidic solution, such as aqueous sulfuric acid, or gaseous carbon dioxide, to obtain the free acids. Composed primarily of tar acids such as phenol, cresols, and xylenols.]	266-019-3	65996-85-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-117-00-1	Tar acids, brown-coal, crude; Crude Phenols; [An acidified alkaline extract of brown coal tar distillate. Composed primarily of phenol and phenol homologs.]	309-888-7	101316-86-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-118-00-7	Tar acids, brown-coal gasification; Crude Phenols; [A complex combination of organic compounds obtained from brown coal gasification. Composed primarily of C_{6-10} hydroxy aromatic phenols and their homologs.]	295-536-7	92062-22-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-119-00-2	Tar acids, distn. residues; Distillate Phenols; [A residue from the distillation of crude phenol from coal. It consists predominantly of phenols having carbon numbers in the range of C_8 through C_{10} with a softening point of 60 °C to 80 °C (140 °F to 176 °F).]	306-251-5	96690-55-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-120-00-8	Tar acids, methylphenol fraction; Distillate Phenols; [The fraction of tar acid rich in 3- and 4-methylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-892-9	84989-04-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-121-00-3	Tar acids, polyalkylphenol fraction; Distillate Phenols; [The fraction of tar aicds, recovered by distillation of low-temperature coal tar crude tar acids, having an approximate boiling range of 225 °C to 320 °C (437 °F to 608 °F). Composed primarily of polyalkylphenols.]	284-893-4	84989-05-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-122-00-9	Tar acids, xylenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 2,4- and 2,5-dimethylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-895-5	84989-06-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-123-00-4	Tar acids, ethylphenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 3- and 4-ethylphenol, recovered by distillation of low-temperature coal tar crude tar acids.]	284-891-3	84989-03-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-124-00-X	Tar acids, 3,5-xylenol fraction; Distillate Phenols; [The fraction of tar acids, rich in 3,5-dimethylphenol, recovered by distillation of low-temperature coal tar acids.]	284-896-0	84989-07-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-125-00-5	Tar acids, residues, distillates, first-cut; Distillate Phenols; [The residue from the distillation in the range of 235 °C to 355 °C (481 °F to 697 °F) of light carbolic oil.]	270-713-1	68477-23-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-126-00-0	Tar acids, cresylic, residues; Distillate Phenols; [The residue from crude coal tar acids after removal of phenol, cresols, xylenols and any higher boiling phenols. A black solid with a melting point approximately 80 °C (176 °F). Composed primarily of polyalkyphenols, resin gums, and inorganic salts.]	271-418-0	68555-24-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-127-00-6	Phenols, C_{9-11}; Distillate Phenols	293-435-2	91079-47-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-128-00-1	Tar acids, cresylic; Distillate Phenols; [A complex combination of organic compounds obtained from brown coal and boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). It contains chiefly phenols and pyridine bases.]	295-540-9	92062-26-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-129-00-7	Tar acids, brown-coal, C_2-alkylphenol fraction; Distillate Phenols; [The distillate from the acidification of alkaline washed lignite tar distillate boiling in the range of approximately 200 °C to 230 °C (392 °F to 446 °F). Composed primarily of <{ITA}>m-<{/ITA}> and <{ITA}>p-<{/ITA}>ethylphenol as well as cresols and xylenols.]	302-662-9	94114-29-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-130-00-2	Extract oils (coal), naphthalene oils; Acid Extract; [The aqueous extract produced by an acidic wash of alkali-washed napthalene oil. Composed primarily of acid salts of various aromatic nitrogen bases including pyridine, quinoline and their alkyl derivatives.]	292-623-1	90641-00-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-131-00-8	Tar bases, quinoline derivs.; Distillate Bases	271-020-7	68513-87-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-132-00-3	Tar bases, coal, quinoline derivs. fraction; Distillate Bases	274-560-1	70321-67-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-133-00-9	Tar bases, coal, distn. residues; Distillate Bases; [The distillation residue remaining after the distillation of the neutralized, acid-extracted base-containing tar fractions obtained by the distillation of coal tars. It contains chiefly aniline, collidines, quinoline and quinoline derivatives and toluidines.]	295-544-0	92062-29-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-134-00-4	Hydrocarbon oils, arom., mixed with polyethylene and polypropylene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of a polyethylene/polypropylene mixture with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of approximately 70 °C to 120 °C (158 °F to 248 °F).]	309-745-9	100801-63-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-135-00-X	Hydrocarbon oils, arom., mixed with polyethylene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of polyethylene with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of 70 °C to 120 °C (158 °F to 248 °F).]	309-748-5	100801-65-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-136-00-5	Hydrocarbon oils, arom., mixed with polystyrene, pyrolyzed, light oil fraction; Heat Treatment Products; [The oil obtained from the heat treatment of polystyrene with coal tar pitch or aromatic oils. It consists predominantly of benzene and its homologs boiling in a range of approximately 70 °C to 210 °C (158 °F to 410 °F).]	309-749-0	100801-66-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-137-00-0	Extract residues (coal), tar oil alk., naphthalene distn. residues; Naphthalene Oil Extract Residue; [The residue obtained from chemical oil extracted after the removal of naphthalene by distillation composed primarily of two to four membered condensed ring aromatic hydrocarbons and aromatic nitrogen bases.]	277-567-8	73665-18-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-138-00-6	Creosote oil, low-boiling distillate; Wash Oil; [The low-boiling distillation fraction obtained from the high temperature carbonization of bituminous coal, which is further refined to remove excess crystalline salts. It consists primarily of creosote oil with some of the normal polynuclear aromatic salts, which are components of coal tar distillate, removed. It is crystal free at approximately 38 °C (100 °F).]	274-566-4	70321-80-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-139-00-1	Tar acids, cresylic, sodium salts, caustic solns.; Alkaline Extract	272-361-4	68815-21-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-140-00-7	Extract oils (coal), tar base; Acid Extract; [The extract from coal tar oil alkaline extract residue produced by an acidic wash such as aqueous sulfuric acid after distillatin to remove naphthalene. Composed primarily of the acid salts of various aromatic nitrogen bases including pyridine, quinoline, and their alkyl derivatives.]	266-020-9	65996-86-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-141-00-2	Tar bases, coal, crude; Crude Tar Bases; [The reaction product obtained by neutralizing coal tar base extract oil with an alkaline solution, such as aqueous sodium hydroxide, to obtain the free bases. Composed primarily of such organic bases as acridine, phenanthridine, pyridine, quinoline and their alkyl derivatives.]	266-018-8	65996-84-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		HJM
648-142-00-8	Residues (coal), liq. solvent extn.; [A cohesive powder composed of coal mineral matter and undissolved coal remaining after extraction of coal by a liquid solvent.]	302-681-2	94114-46-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-143-00-3	Coal liquids, liq. solvent extn. soln.; [The product obtained by filtration of coal mineral matter and undissolved coal from coal extract solution produced by digesting coal in a liquid solvent. A black, viscous, highly complex liquid combination composed primarily of aromatic and partly hydro-genated aromatic hydrocarbons, aromatic nitrogen compounds, aromatic sulfur compounds, phenolic and other aromatic oxygen compounds and their alkyl derivatives.]	302-682-8	94114-47-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M
648-144-00-9	Coal liquids, liq. solvent extn.; [The substantially solvent-free product obtained by the distillation of the solvent from filtered coal extract solution produced by digesting coal in a liquid solvent. A black semi-solid, composed primarily of a complex combination of condensed-ring aromatic hydrocarbons, aromatic nitrogen compounds, aromatic sulfur compounds, phenolic compounds and other aromatic oxygen compounds, and their alkyl derivatives.]	302-683-3	94114-48-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H M

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-145-00-4	Tar brown-coal; [An oil distilled from brown-coal tar. Composed primarily of aliphatic, naphthenic and one- to three-ring aromatic hydrocarbons, their alkyl derivates, heteroaromatics and one- and two-ring phenols boiling in the range of approximately 150 °C to 360 °C (302 °F to 680 °F).]	309-885-0	101316-83-0	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
648-146-00-X	Tar, brown-coal, low-temp.; [A tar obtained from low temperature carbonization and low temperature gasification of brown coal. Composed primarily of aliphatic, naphthenic and cyclic aromatic hydrocarbons, heteroaromatic hydrocarbons and cyclic phenols.]	309-886-6	101316-84-1	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
648-147-00-5	Light oil (coal), coke-oven; Crude benzole; [The volatile organic liquid extracted from the gas evolved in the high temperature (greater than 700 °C (1292 °F)) destructive distillation of coal. Composed primarily of benzene, toluene, and xylenes. May contain other minor hydrocarbon constituents.]	266-012-5	65996-78-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-148-00-0	Distillates (coal), liq. solvent extn., primary; [The liquid product of condensation of vapors emitted during the digestion of coal in a liquid solvent and boiling in the range of approximately 30 °C to 300 °C (86 °F to 572 °F). Composed primarily of partly hydrogenated condensed-ring aromatic hydrocarbons, aromatic compounds containing nitrogen, oxygen and sulfur, and their alkyl derivatives having carbon numbers predominantly in the range of C_4 through C_{14}.]	302-688-0	94114-52-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-149-00-6	Distillates (coal), solvent extn., hydrocracked; [Distillate obtained by hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction process and boiling in the range of approximately 30 °C to 300 °C (86 °F to 572 °F). Composed primarily of aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes with carbon numbers predominantly in the range of C_4 through C_{14}. Nitrogen, sulfur and oxygen-containing aromatic and hydrogenated aromatic compounds are also present.]	302-689-6	94114-53-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-150-00-1	Naphtha (coal), solvent extn., hydrocracked; [Fraction of the distillate obtained by hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes with carbon numbers predominantly in the range of C_4 to C_9. Nitrogen, sulfur and oxygen-containing aromatic and hydrogenated aromatic compounds are also present.]	302-690-1	94114-54-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-151-00-7	Gasoline, coal solvent extn., hydrocracked naphtha; [Motor fuel produced by the reforming of the refined naphtha fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 30 °C to 180 °C (86 °F to 356 °F). Composed primarily of aromatic and naphthenic hydrocarbons, their alkyl derivatives and alkyl hydrocarbons having carbon numbers in the range of C_4 through C_9.]	302-691-7	94114-55-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
648-152-00-2	Distillates (coal), solvent extn., hydrocracked middle; [Distillate obtained from the hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 300 °C (356 °F to 572 °F). Composed primarily of two-ring aromatic, hydrogenated aromatic and naphthenic compounds, their alkyl derivatives and alkanes having carbon numbers predominantly in the range of C_9 through C_{14}. Nitrogen, sulfur and oxygen-containing compounds are also present.]	302-692-2	94114-56-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
648-153-00-8	Distillates (coal), solvent extn., hydrocracked hydrogenated middle; [Distillate from the hydrogenation of hydrocracked middle distillate from coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 280 °C (356 °F to 536 °F). Composed primarily of hydrogenated two- ring carbon compounds and their alkyl derivatives having carbon numbers predominantly in the range of C_9 through C_{14}.]	302-693-8	94114-57-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
648-154-00-3	Fuels, jet aircraft, coal solvent extn., hydrocracked hydrogenated; [Jet engine fuel produced by hydrogenation of the middle distillate fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 180 °C to 225 °C (356 °F to 473 °F). Composed primarily of hydrogenated two-ring hydrocarbons and their alkyl derivatives having carbon numbers predominantly in the range of C_{10} through C_{12}.]	302-694-3	94114-58-6	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
648-155-00-9	Fuels, diesel, coal solvent extn., hydrocracked hydrogenated; [Diesel engine fuel produced by the hydrogenation of the middle distillate fraction of the products of hydrocracking of coal extract or solution produced by the liquid solvent extraction or supercritical gas extraction processes and boiling in the range of approximately 200 °C to 280 °C (392 °F to 536 °F). Composed primarily of hydrogenated two-ring hydrocarbons and their alkyl derivatives having carbon numbers predominantly in the range of C_{11} through C_{14}.]	302-695-9	94114-59-7	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
648-156-00-4	Light oil (coal), semi-coking process; Fresh oil; [The volatile organic liquid condensed from the gas evolved in the low temperature (less than 700 °C (1292 °F) destructive distillation of coal. Composed primarily of C_{6-10} hydrocarbons.]	292-635-7	90641-11-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H J
649-001-00-3	Extracts (petroleum), light naphthenic distillate solvent	265-102-1	64742-03-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-002-00-9	Extracts (petroleum), heavy paraffinic distillate solvent	265-103-7	64742-04-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-003-00-4	Extracts (petroleum), light paraffinic distillate solvent	265-104-2	64742-05-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-004-00-X	Extracts (petroleum), heavy naphthenic distillate solvent	265-111-0	64742-11-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-005-00-5	Extracts (petroleum), light vacuum gas oil solvent	295-341-7	91995-78-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-006-00-0	hydrocarbons C_{26-55}, arom-rich	307-753-7	97722-04-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-007-00-6	fatty acids, tall-oil, reaction products with iminodiethanol and boric acid	400-160-5	—	Xi; R38 N; R51-53	Xi; N R: 38-51/53 S: (2-)28-37-61		
649-008-00-1	Residues (petroleum), atm. tower; Heavy Fuel oil; [A complex residuum from the atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-045-2	64741-45-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-009-00-7	Gas oils (petroleum), heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and boiling in the range of approximately 350 °C to 600 °C (662 °F to 1112 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	265-058-3	64741-57-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-010-00-2	Distillates (petroleum), heavy catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{35} and boiling in the range of approximately 260 °C to 500 °C (500 °F to 932 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-063-0	64741-61-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-011-00-8	Clarified oils (petroleum), catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-064-6	64741-62-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-012-00-3	Residues (petroleum), hydrocracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the products of a hydrocracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F).]	265-076-1	64741-75-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-013-00-9	Residues (petroleum), thermal cracked; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the product from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-081-9	64741-80-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-014-00-4	Distillates (petroleum), heavy thermal cracked; Heavy Fuel oil; [A complex combination of hydrocarbons from the distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{36} and boiling in the range of approximately 260 °C to 480 °C (500 °F to 896 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-082-4	64741-81-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-015-00-X	Gas oils (petroleum), hydrotreated vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{50} and boiling in the range of approximately 230 °C to 600 °C (446 °F to 1112 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-162-9	64742-59-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-016-00-5	Residues (petroleum), hydrodesulfurized atmospheric tower; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating an atmospheric tower residuum with hydrogen in the presence of a catalyst under conditions primarily to remove organic sulfur compounds. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-181-2	64742-78-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-017-00-0	Gas oils (petroleum), hydrodesulfurized heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and boiling in the range of approximately 350 °C to 600 °C (662 °F to 1112 °C). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-189-6	64742-86-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-018-00-6	Residues (petroleum), steam-cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained as the residual fraction from the distillation of the products of a steam cracking process (including steam cracking to produce ethylene). It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly greater than C_{14} and boiling above approximately 260 °C (500 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	265-193-8	64742-90-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-019-00-1	Residues (petroleum), atmospheric; Heavy Fuel oil; [A complex residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{11} and boiling above approximately 200 °C (392 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	269-777-3	68333-22-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-020-00-7	Clarified oils (petroleum), hydrodesulfurized catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating catalytic cracked clarified oil with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt. % or more of 4-to 6-membered condensed ring aromatic hydrocarbons.]	269-782-0	68333-26-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-021-00-2	Distillates (petroleum), hydrodesulfurized intermediate catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating intermediate catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{30} and boiling in the range of approximately 205 °C to 450 °C (401 °F to 842 °F). It contains a relatively large proportion of tricyclic aromatic hydrocarbons.]	269-783-6	68333-27-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-022-00-8	Distillates (petroleum), hydrodesulfurized heavy catalytic cracked; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treatment of heavy catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{35} and boiling in the range of approximately 260 °C to 500 °C (500 °F to 932 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	269-784-1	68333-28-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-023-00-3	Fuel oil, residues-straight-run gas oils, high-sulfur; Heavy Fuel oil	270-674-0	68476-32-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-024-00-9	Fuel oil, residual; Heavy Fuel oil; [The liquid product from various refinery streams, usually residues. The composition is complex and varies with the source of the crude oil.]	270-675-6	68476-33-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-025-00-4	Residues (petroleum), catalytic reformer fractionator residue distn.; Heavy Fuel oil; [A complex residuum from the distillation of catalytic reformer fractionator residue. It boils approximately above 399 °C (750 °F).]	270-792-2	68478-13-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-026-00-X	Residues (petroleum), heavy coker gas oil and vacuum gas oil; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of heavy coker gas oil and vacuum gas oil. It predominantly consists of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-796-4	68478-17-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-027-00-5	Residues (petroleum), heavy coker and light vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of heavy coker gas oil and light vacuum gas oil. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-983-0	68512-61-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-028-00-0	Residues (petroleum), light vacuum; Heavy Fuel oil; [A complex residuum from the vacuum distillation of the residuum from the atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly greater than C_{13} and boiling above approximately 230 °C (446 °F).]	270-984-6	68512-62-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-029-00-6	Residues (petroleum), steam-cracked light; Heavy Fuel oil; [A complex residuum from the distillation of the products from a steam-cracking process. It consists predominantly of aromatic and unsaturated hydrocarbons having carbon numbers greater than C_7 and boiling in the range of approximately 101 °C to 555 °C (214 °F to 1030 °F).]	271-013-9	68513-69-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-030-00-1	Fuel oil, No 6; Heavy Fuel oil; [A distillate oil having a minimum viscosity of 900 SUS at 37.7 °C (100 °F) to a maximum of 9000 SUS at 37.7 °C (100 °F).]	271-384-7	68553-00-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-031-00-7	Residues (petroleum), topping plant, low-sulfur; Heavy Fuel oil; [A low-sulfur complex combination of hydrocarbons produced as the residual fraction from the topping plant distillation of crude oil. It is the residuum after the straight-run gasoline cut, kerosene cut and gas oil cut have been removed.]	271-763-7	68607-30-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-032-00-2	Gas oils (petroleum), heavy atmospheric; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{35} and boiling in the range of approximately 121 °C to 510 °C (250 °F to 950 °F).]	272-184-2	68783-08-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-033-00-8	Residues (petroleum), coker scrubber, Condensed-ring-arom.-contg.; Heavy Fuel oil; [A very complex combination of hydrocarbons produced as the residual fraction from the distillation of vaccum residuum and the products from a thermal cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{20} and boiling above approximately 350 °C (662 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed rind aromatic hydrocarbons.]	272-187-9	68783-13-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-034-00-3	Distillates (petroleum), petroleum residues vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from the atmospheric distillation of crude oil.]	273-263-4	68955-27-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-035-00-9	Residues (petroleum), steam-cracked, resinous; Heavy Fuel oil; [A complex residuum from the distillation of steam-cracked petroleum residues.]	273-272-3	68955-36-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-036-00-4	Distillates (petroleum), intermediate vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum, distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{42} and boiling in the range of approximately 250 °C to 545 °C (482 °F to 1013 °F). This stream is likely to contain 5 wt. % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	274-683-0	70592-76-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-037-00-X	Distillates (petroleum), light vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{35} and boiling in the range of approximately 250 °C to 545 °C (482 °F to 1013 °F).]	274-684-6	70592-77-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-038-00-5	Distillates (petroleum), vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having numbers predominantly in the range of C_{15} through C_{50} and boiling in the range of approximately 270 °C to 600 °C (518 °F to 1112 °F). This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	274-685-1	70592-78-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-039-00-0	Gas oils (petroleum), hydrodesulfurized coker heavy vacuum; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by hydrodesulfurization of heavy coker distillate stocks, It consists predominantly of hydrocarbons having carbon numbers predominantly in the range C_{18} to C_{44} and boiling in the range of approximately 304 °C to 548 °C (579 °F to 1018 °F). Likely to contain 5 % or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	285-555-9	85117-03-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-040-00-6	Residues (petroleum), steam-cracked, distillates; Heavy Fuel oil; [A complex combination of hydrocarbons obtained during the production of refined petroleum tar by the distillation of steam cracked tar. It consists predominantly of aromatic and other hydrocarbons and organic sulfur compounds.]	292-657-7	90669-75-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-041-00-1	Residues (petroleum), vacuum, light; Heavy Fuel oil; [A complex residuum from the vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{24} and boiling above approximately 390 °C (734 °F).]	292-658-2	90669-76-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-042-00-7	Fuel oil, heavy, high-sulfur; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the distillation of crude petroleum. It consists predominantly of aliphatic, aromatic and cycloaliphatic hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	295-396-7	92045-14-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-043-00-2	Residues (petroleum), catalytic cracking; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from the distillation of the products from a catalytic cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{11} and boiling above approximately 200 °C (392 °F).]	295-511-0	92061-97-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-044-00-8	Distillates (petroleum), intermediate catalytic cracked, thermally degraded; Heavy Fuel oil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process which has been used as a heat transfer fluid. It consists predominantly of hydrocarbons boiling in the range of approximately 220 °C to 450 °C (428 °F to 842 °F). This stream is likely to contain organic sulfur compounds.]	295-990-6	92201-59-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-045-00-3	Residual oils (petroleum); Heavy Fuel oil; [A complex combination of hydrocarbons, sulfur compounds and metal-containing organic compounds obtained as the residue from refinery fractionation cracking processes. It produces a finished oil with a viscosity above 2cSt. at 100 °C.]	298-754-0	93821-66-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-046-00-9	Residues, steam cracked, thermally treated; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by the treatment and distillation of raw steam-cracked naphtha. It consists predominantly of unsaturated hydrocarbons boiling in the range above approximately 180 °C (356 °F).]	308-733-0	98219-64-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-047-00-4	Distillates (petroleum), hydrodesulfurized full-range middle; Heavy Fuel oil; [A complex combination of hydrocarbons obtained by treating a petroleum stock with hydrogen. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F).]	309-863-0	101316-57-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-048-00-X	Residues (petroleum), catalytic reformer fractionator; Heavy Fuel oil; [A complex combination of hydrocarbons produced as the residual fraction from distillation of the product from a catalytic reforming process. It consists of predominantly aromatic hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{25} and boiling in the range of approximately 160 °C to 400 °C (320 °F to 725 °F). This stream is likely to contain 5 wt. % or more of 4- or 6-membered condensed ring aromatic hydrocarbons.]	265-069-3	64741-67-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-049-00-5	Petroleum; Crude oil; [A complex combination of hydrocarbons, It consists predominantly of aliphatic, alicyclic and aromatic hydrocarbons. It may also contain small amounts of nitrogen, oxygen and sulfur compounds. This category encompasses light, medium, and heavy petroleums, as well as the oils extended from tar sands. Hydrocarbonaceous materials requiring major chemical changes for their recovery or conversion to petroleum refinery feedstocks such as crude shale oils; upgraded shale oils and liquid coal fuels are not included in this definition.]	232-298-5	8002-05-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-050-00-0	Distillates (petroleum), light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated aliphatic hydrocarbons normally present in this distillation range of crude oil.]	265-051-5	64741-50-0	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-051-00-6	Distillates (petroleum), heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated aliphatic hydrocarbons.]	265-052-0	64741-51-1	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-052-00-1	Distillates (petroleum), light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-053-6	64741-52-2	Carc. Cat. 1; R45	T R: 45 S: 53-45		H

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649-053-00-7	Distillates (petroleum), heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by vacuum distillation of the residuum from atmospheric distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-054-1	64741-53-3	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-054-00-2	Distillates (petroleum), acid-treated heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-117-3	64742-18-3	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-055-00-8	Distillates (petroleum), acid-treated light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-118-9	64742-19-4	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-056-00-3	Distillates (petroleum), acid-treated heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of a least 100 SUS at 100 °F (19cSt at 40 °C).]	265-119-4	64742-20-7	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-057-00-9	Distillates (petroleum), acid-treated light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-121-5	64742-21-8	Carc. Cat. 1; R45	T R: 45 S: 53-45		H

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649-058-00-4	Distillates (petroleum), chemically neutralized heavy paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons obtained from a treating process to remove acidic materials. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of aliphatic hydrocarbons.]	265-127-8	64742-27-4	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-059-00-X	Distillates (petroleum), chemically neutralized light paraffinic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-128-3	64742-28-5	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-060-00-5	Distillates (petroleum), chemically neutralized heavy naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-135-1	64742-34-3	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-061-00-0	Distillates (petroleum), chemically neutralized light naphthenic; Unrefined or mildly refined baseoil; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS a 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-136-7	64742-35-4	Carc. Cat. 1; R45	T R: 45 S: 53-45		H
649-062-00-6	Gases (petroleum), catalytic cracked naphtha depropanizer overhead, C_3-rich acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked hydrocarbons and treated to remove acidic impurities. It consists of hydrocarbons having carbon numbers in the range of C_2 through C_4, predominantly C_3.]	270-755-0	68477-73-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-063-00-1	Gases (petroleum), catalytic cracker; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic cracking process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-756-6	68477-74-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-064-00-7	Gases (petroleum), catalytic cracker, C_{1-5}-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_6, predominantly C_1 through C_5.]	270-757-1	68477-75-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-065-00-2	Gases (petroleum), catalytic polymd. naphtha stabilizer overhead, C_{2-4}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of catalytic polymerized naphtha. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_2 through C_6, predominantly C_2 through C_4.]	270-758-7	68477-76-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-066-00-8	Gases (petroleum), catalytic reformer, C_{1-4}-rich; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers in the range of C_1 through C_6, predominantly C_1 through C_4.]	270-760-8	68477-79-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-067-00-3	Gases (petroleum), C_{3-5} olefinic-paraffinic alkylation feed; Petroleum gas; [A complex combination of olefinic and paraffinic hydrocarbons having carbon numbers in the range of C_3 through C_5 which are used as alkylation feed. Ambient temperatures normally exceed the critical temperature of these combinations.]	270-765-5	68477-83-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-068-00-9	Gases (petroleum), C_4-rich; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from a catalytic fractionation process. It consists of aliphatic hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_4.]	270-767-6	68477-85-0	⊗ ⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-069-00-4	Gases (petroleum), deethanizer overheads; Petroleum gas; [A complex combination of hydrocarbons produced from distillation of the gas and gasoline fractions from the catalytic cracking process. It contains predominantly ethane and ethylene.]	270-768-1	68477-86-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-070-00-X	Gases (petroleum), deisobutanizer tower overheads; Petroleum gas; [A complex combination of hydrocarbons produced by the atmospheric distillation of a butane-butylene stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	270-769-7	68477-87-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-071-00-5	Gases (petroleum), depropanizer dry, propene-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from the gas and gasoline fractions of a catalytic cracking process. It consists predominantly of propylene with some ethane and propane.]	270-772-3	68477-90-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-072-00-0	Gases (petroleum), depropanizer overheads; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from the gas and gasoline fractions of a catalytic cracking process. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	270-773-9	68477-91-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-073-00-6	Gases (petroleum), gas recovery plant depropanizer overheads; Petroleum gas; [A complex combination of hydrocarbons obtained by fractionation of miscellaneous hydrocarbon streams. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4, predominantly propane.]	270-777-0	68477-94-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-074-00-1	Gases (petroleum), Girbatol unit feed; Petroleum gas; [A complex combination of hydrocarbons that is used as the feed into the Girbatol unit to remove hydrogen sulfide. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	270-778-6	68477-95-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-075-00-7	Gases (petroleum), isomerized naphtha fractionator, C_4-rich, hydrogen sulfide-free; Petroleum gas	270-782-8	68477-99-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-076-00-2	Tail gas (petroleum), catalytic cracked clarified oil and thermal cracked vacuum residue fractionation reflux drum; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked clarified oil and thermal cracked vacuum residue. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-802-5	68478-21-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-077-00-8	Tail gas (petroleum), catalytic cracked naphtha stabilization absorber; Petroleum gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-803-0	68478-22-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-078-00-3	Tail gas (petroleum), catalytic cracker, catalytic reformer and hydrodesulfurizer combined fractionater; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of products from catalytic cracking, catalytic reforming and hydrodesulfurizing processes treated to remove acidic impurities. It consists predominantly of hydrocarbons having cabon numbers predominantly in the range of C_1 through C_5.]	270-804-6	68478-24-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-079-00-9	Tail gas (petroleum), catalytic reformed naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of catalytic reformed naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	270-806-7	68478-26-2	⊗Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-080-00-4	Tail gas (petroleum), saturate gas plant mixed stream, C_4-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of straight-run naphtha, distillation tail gas and catalytic reformed naphtha stabilizer tail gas. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly butane and isobutane.]	270-813-5	68478-32-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-081-00-X	Tail gas (petroleum), saturate gas recovery plant, C_{1-2}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of distillate tail gas, straight-run naphtha, catalytic reformed naphtha stabilizer tail gas. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1through C_5, predominantly methane and ethane.]	270-814-0	68478-33-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-082-00-5	Tail gas (petroleum), vacuum residues thermal cracker; Petroleum gas; [A complex combination of hydrocarbons obtained from the thermal cracking of vacuum residues. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-815-6	68478-34-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-083-00-0	Hydrocarbons, C_{3-4}-rich, petroleum distillate; Petroleum gas; [A complex combination of hydrocarbons produced by distillation and condensation of crude oil. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_3 through C_4.]	270-990-9	68512-91-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-084-00-6	Gases (petroleum), full-range straight-run naphtha dehexanizer off; petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of the full-range straight-run naphtha. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-000-8	68513-15-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-085-00-1	Gases (petroleum), hydrocracking depropanizer off, hydrocarbon-rich; Petroleum gas; [A complex combination of hydrocarbon produced by the distillation of products from a hydrocracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4. It may also contain small amounts of hydrogen and hydrogen sulfide.]	271-001-3	68513-16-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-086-00-7	Gases (petroleum), light straight-run naphtha stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the stabilization of light straight-run naphtha. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-002-9	68513-17-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-087-00-2	Residues (petroleum), alkylation splitter, C_4-rich; Petroleum gas; [A complex residuum from the distillation of streams various refinery operations. It consists of hydrocarbons having carbon numbers in the range of C_4 through C_5, predominantly butane and boiling in the range of approximately - 11.7 °C to 27.8 °C (11 °F to 82 °F).]	271-010-2	68513-66-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-088-00-8	Hydrocarbons, C_{1-4}; Petroleum gas; [A complex combination of hydrocarbons provided by thermal cracking and absorber operations and by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately minus 164 °C to minus 0.5 °C (-263 °F to 31 °F).]	271-032-2	68514-31-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-089-00-3	Hydrocarbons, C_{1-4}, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting hydrocarbon gases to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately - 164 °C to - 0.5 °C (-263 °F to 31 °F).]	271-038-5	68514-36-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-090-00-9	Hydrocarbons, C_{1-3}; Petroleum gas; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3 and boiling in the range of approximately minus 164 °C to minus 42 °C (-263 °F to - 44 °F).]	271-259-7	68527-16-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-091-00-4	Hydrocarbons, C_{1-4}, debutanizer fraction; Petroleum gas	271-261-8	68527-19-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-092-00-X	Gases (petroleum), C_{1-5}, wet; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of crude oil and/or the cracking of tower gas oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	271-624-0	68602-83-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-093-00-5	Hydrocarbons, C_{2-4}; Petroleum gas	271-734-9	68606-25-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-094-00-0	Hydrocarbons, C_3; Petroleum gas	271-735-4	68606-26-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-095-00-6	Gases (petroleum), alkylation feed; Petroleum gas; [A complex combination of hydrocarbons produced by the catalytic cracking of gas oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	271-737-5	68606-27-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-096-00-1	Gases (petroleum), depropanizer bottoms fractionation off; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of depropanizer bottoms. It consists predominantly of butane, isobutane and butadiene.]	271-742-2	68606-34-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-097-00-7	Gases (petroleum), refinery blend; Petroleum gas; [A complex combination obtained from various processes. It consists of hydrogen, hydrogen sulfide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-183-7	68783-07-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-098-00-2	Gases (petroleum), catalytic cracking; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5.]	272-203-4	68783-64-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-099-00-8	Gases (petroleum), C_{2-4}, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4 and boiling in the range of approximately - 51 °C to - 34 °C (-60 °F to - 30 °F).]	272-205-5	68783-65-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-100-00-1	Gases (petroleum), crude oil fractionation off; Petroleum gas; [A complex combination of hydrocarbons produced by the fractionation of crude oil. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-871-7	68918-99-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-101-00-7	Gases (petroleum), dehexanizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of combined naphtha streams. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-872-2	68919-00-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-102-00-2	Gases (petroleum), light straight run gasoline fractionation stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of light straight-run gasoline. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-878-5	68919-05-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-103-00-8	Gases (petroleum), naphtha unifiner desulfurization stripper off; Petroleum gas; [A complex combination of hydrocarbons produced by a naphtha unifiner desulfurization process and stripped from the naphtha product. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-879-0	68919-06-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-104-00-3	Gases (petroleum), straight-run naphtha catalytic reforming off; Petroleum gas; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha and fractionation of the total effluent. It consists of methane, ethane, and propane.]	272-882-7	68919-09-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-105-00-9	Gases (petroleum), fluidized catalytic cracker splitter over-heads; Petroleum gas; [A complex combination of hydrocarbons produced by the fractionation of the charge to the C_3 -C_4 splitter. It consists predominantly of C_3 hydrocarbons.]	272-893-7	68919-20-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-106-00-4	Gases (petroleum), straight-run stabilizer off; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation of the liquid from the first tower used in the distillation of crude oil. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-883-2	68919-10-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45 S: 53-45		H K

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649-107-00-X	Gases (petroleum), catalytic cracked naphtha debutanizer; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked naphtha. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-169-3	68952-76-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-108-00-5	Tail gas (petroleum), catalytic cracked distillate and naphtha stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained by the fractionation of catalytic cracked naphtha and distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-170-9	68952-77-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-109-00-0	Tail gas (petroleum), thermal-cracked distillate, gas oil and naphtha absorber; petroleum gas; [A complex combination of hydrocarbons obtained from the separation of thermal-cracked distillates, naphtha and gas oil. It consists pedrominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-175-6	68952-81-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-110-00-6	Tail gas (petroleum), thermal cracked hydrocarbon fractionation stabilizer, petroleum coking; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization of thermal cracked hydrocarbons from petroleum coking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-176-1	68952-82-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-111-00-1	Gases (petroleum, light steam-cracked, butadiene conc.; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a thermal cracking process, It consists of hydrocarbons having a carbon number predominantly of C_4.]	273-265-5	68955-28-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-112-00-7	Gases (petroleum), straight-run naphtha catalytic reformer stabilizer overhead; Petroleum gas; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha and the fractionation of the total effluent. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	273-270-2	68955-34-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-113-00-2	Hydrocarbons, C_4; Petroleum gas	289-339-5	87741-01-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-114-00-8	Alkanes, C_{1-4}, C_3-rich; Petroleum gas	292-456-4	90622-55-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-115-00-3	Gases (petroleum), steam-cracker C_3-rich; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a steam cracking process. It consists predominantly of propylene with some propane and boils in the range of approximately - 70 °C to 0 °C (-94 °F to 32 °F).]	295-404-9	92045-22-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-116-00-9	Hydrocarbons, C_4, steam-cracker distillate; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of the products of a steam cracking process. It consists predominantly of hydrocarbons having a carbon number of C_4, predominantly 1-butene and 2-butene, containing also butane and isobutene and boiling in the range of approximately minus 12 °C to 5 °C (10.4 °F to 41 °F).]	295-405-4	92045-23-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-117-00-4	Petroleum gases, liquefied, sweetened, C_4 fraction; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting a liquified petroleum gas mix to a sweetening process to oxidize mercaptans or to remove acidic impurities. It consists predominantly of C_4 saturated and unsaturated hydrocarbons.]	295-463-0	92045-80-2	F+; R12 Carc. Cat. 1; R45 Muta. Cat. 2; R46	F+; T R: 12-45-46 S: 53-45		HKS
649-118-00-X	Hydrocarbons, C_4, 1,3-butadiene- and isobutene-free; Petroleum gas	306-004-1	95465-89-7	⊗ Carc. Cat. 2; R45	T R: 45 S: 53-45		H K
649-119-00-5	Raffinates (petroleum), steam-cracked C_4 fraction cuprous ammonium acetate extn., C_{3-5} and C_{3-5} unsatd., butadiene-free; Petroleum gas	307-769-4	97722-19-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-120-00-0	Gases (petroleum), amine system feed; Refinery gas; [The feed gas to the amine system for removal of hydrogen sulfide. It consists of hydrogen. Carbon monoxide, carbon dioxide, hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5 may also be present.]	270-746-1	68477-65-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-121-00-6	Gases (petroleum), benzene unit hydrodesulfurizer off; Refinery gas; [Off gases produced by the benzene unit. It consists primarily of hydrogen. Carbon monoxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6, including benzene, may also be present.]	270-747-7	68477-66-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-122-00-1	Gases (petroleum), benzene unit recycle, hydrogen-rich; Refinery gas; [A complex combination of hydrocarbons obtained by recycling the gases of the benzene unit. It consists primarily of hydrogen with various small amounts of carbon monoxide and hydrocarbons having carbon numbers in the range of C_1 through C_6.]	270-748-2	68477-67-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-123-00-7	Gases (petroleum), blend oil, hydrogen-nitrogen-rich; Refinery gas; [A complex combination of hydrocarbons obtained by distillation of a blend oil. It consists primarily of hydrogen and nitrogen with various small amounts of carbon monoxide, carbon dioxide, and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-749-8	68477-68-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-124-00-2	Gases (petroleum), catalytic reformed naphtha stripper overheads; Refinery gas; [A complex combination of hydrocarbons obtained from stabilization of catalytic reformed naphtha. Its consists of hydrogen and saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	270-759-2	68477-77-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-125-00-8	Gases (petroleum), C_{6-8} catalytic reformer recycle; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from catalytic reforming of C_6-C_8 feed and recycled to conserve hydrogen. It consists primarily of hydrogen. It may also contain various small amounts of carbon monoxide, carbon dioxide, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-761-3	68477-80-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-126-00-3	Gases (petroleum), C_{6-8} catalytic reformer; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from catalytic reforming of C_6-C_8 feed. It consists of hydrocarbons having carbon numbers in the range of C_1 through C_5 and hydrogen.]	270-762-9	68477-81-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-127-00-9	Gases (petroleum), C_{6-8} catalytic reformer recycle, hydrogen-rich; Refinery gas	270-763-4	68477-82-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-128-00-4	Gases (petroleum), C_2-return stream; Refinery gas; [A complex combination of hydrocarbons obtained by the extraction of hydrogen from a gas stream which consists primarily of hydrogen with small amounts of nitrogen, carbon monoxide, methane, ethane, and ethylene. It contains predominantly hydrocarbons such as methane, ethane, and ethylene with small amounts of hydrogen, nitrogen and carbon monoxide.]	270-766-0	68477-84-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-129-00-X	Gases (petroleum), dry sour, gas-concn.-unit-off; Refinery gas; [The complex combination of dry gases from a gas concentration unit. It consists of hydrogen, hydrogen sulfide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	270-774-4	68477-92-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-130-00-5	Gases (petroleum), gas concn. reabsorber distn.; Refinery gas; [A complex combination of hydrocarbons produced by distillation of products from combined gas streams in a gas concentration reabsorber. It consists predominantly of hydrogen, carbon monoxide, carbon dioxide, nitrogen, hydrogen sulfide and hydrocarbons having carbon numbers in the range of C_1 through C_3.]	270-776-5	68477-93-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-131-00-0	Gases (petroleum), hydrogen absorber off; Refinery gas; [A complex combination obtained by absorbing hydrogen from a hydrogen rich stream. It consists of hydrogen, carbon monoxide, nitrogen, and methane with small amounts of C_2 hydrocarbons.]	270-779-1	68477-96-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-132-00-6	Gases (petroleum), hydrogen-rich; Refinery gas; [A complex combination separated as a gas from hydrocarbon gases by chilling. It consists primarily of hydrogen with various small amounts of carbon monoxide, nitrogen, methane, and C_2 hydrocarbons.]	270-780-7	68477-97-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-133-00-1	Gases (petroleum), hydrotreater blend oil recycle, hydrogen-nitrogen-rich; Refinery gas; [A complex combination obtained from recycled hydrotreated blend oil. It consists primarily of hydrogen and nitrogen with various small amounts of carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-781-2	68477-98-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-134-00-7	Gases (petroleum), recycle, hydrogen-rich; Refinery gas; [A complex combination obtained from recycled reactor gases. It consists primarily of hydrogen with various small amounts of carbon monoxide, carbon dioxide, nitrogen, hydrogen sulfide, and saturated aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_5.]	270-783-3	68478-00-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-135-00-2	Gases (petroleum), reformer make-up, hydrogen-rich; Refinery gas; [A complex combination obtained from the reformers. It consists primarily of hydrogen with various small amounts of carbon monoxide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-784-9	68478-01-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-136-00-8	Gases (petroleum), reforming hydrotreater; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen, methane, and ethane with various small amounts of hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 thorugh C_5.]	270-785-4	68478-02-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-137-00-3	Gases (petroleum), reforming hydrotreater, hydrogen-methane-rich; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen and methane with various small amounts of carbon monoxide, carbon dioxide, nitrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_5.]	270-787-5	68478-03-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-138-00-9	Gases (petroleum), reforming hydrotreater make-up, hydrogen-rich; Refinery gas; [A complex combination obtained from the reforming hydrotreating process. It consists primarily of hydrogen with various small amounts of carbon monoxide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-788-0	68478-04-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-139-00-4	Gases (petroleum), thermal cracking distn.; Refinery gas; [A complex combination produced by distillation of products from a thermal cracking process. It consists of hydrogen, hydrogen sulfide, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-789-6	68478-05-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-140-00-X	Tail gas (petroleum), catalytic cracker refractionation absorber; Refinery gas; [A complex combination of hydrocarbons obtained from refractionation of products from a catalytic cracking process. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	270-805-1	68478-25-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-141-00-5	Tail gas (petroleum), catalytic reformed naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from the catalytic reforming of straight run naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-807-2	68478-27-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-142-00-0	Tail gas (petroleum), catalytic reformed naphtha stabilizer; Refinery gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic reformed naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-808-8	68478-28-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-143-00-6	Tail gas (petroleum), cracked distillate hydrotreater separator; Refinery gas; [A complex combination of hydrocarbons obtained by treating cracked distillates with hydrogen in the presence of a catalyst. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	270-809-3	68478-29-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-144-00-1	Tail gas (petroleum), hydrodesulfurized straight-run naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from hydrodesulfurization of straight-run naphtha. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	270-810-9	68478-30-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-145-00-7	Gases (petroleum), catalytic reformed straight-run naphtha stabilizer overheads; Refinery gas; [A complex combination of hydrocarbons obtained from the catalytic reforming of straight-run naphtha followed by fractionation of the total effluent. It consists of hydrogen, methane, ethane and propane.]	270-999-8	68513-14-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-146-00-2	Gases (petroleum), reformer effluent high-pressure flash drum off; Refinery gas; [A complex combination produced by the high-pressure flashing of the effluent from the reforming reactor. It consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-003-4	68513-18-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-147-00-8	Gases (petroleum), reformer effluent low-pressure flash drum off; Refinery gas; [A complex combination produced by low-pressure flashing of the effluent from the reforming reactor. It consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-005-5	68513-19-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-148-00-3	Gases (petroleum), oil refinery gas distn. off; Refinery gas; [A complex combination separated by distillation of a gas stream containing hydrogen, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers in the range of C_1 through C_6 or obtained by cracking ethane and propane. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_2, hydrogen, nitrogen, and carbon monoxide.]	271-258-1	68527-15-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-149-00-9	Gases (petroleum), benzene unit hydrotreater dependentanizer overheads; Refinery gas; [A complex combination produced by treating the feed from the benzene unit with hydrogen in the presence of a catalyst followed by dependentanizing. It consists primarily of hydrogen, ethane and propane with various small amounts of nitrogen, carbon monoxide, carbon dioxide and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6. It may contain trace amounts of benzene.]	271-623-5	68602-82-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-150-00-4	Gases (petroleum), secondary absorber off, fluidized catalytic cracker overheads fractionator; Refinery gas; [A complex combination produced by the fractionation of the overhead products from the catalytic cracking process in the fluidized catalytic cracker. It consists of hydrogen, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	271-625-6	68602-84-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-151-00-X	Petroleum products, refinery gases; Refinery gas; [A complex combination which consists primarily of hydrogen with various small amounts of methane, ethane, and propane.]	271-750-6	68607-11-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-152-00-5	Gases (petroleum), hydrocracking low-pressure separator; Refinery gas; [A complex combination obtained by the liquid-vapor separation of the hydrocracking process reactor effluent. It consists predominantly of hydrogen and saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	272-182-1	68783-06-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-153-00-0	Gases (petroleum), refinery; Refinery gas; [A complex combination obtained from various petroleum refining operations. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	272-338-9	68814-67-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-154-00-6	Gases (petroleum), platformer products separator off; Refinery gas; [A complex combination obtained from the chemical reforming of naphthenes to aromatics. It consists of hydrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_4.]	272-343-6	68814-90-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-155-00-1	Gases (petroleum), hydrotreated sour kerosine depentanizer stabilizer off; Refinery gas; [The complex combination obtained from the depentanizer stabilization of hydrotreated kerosine. It consists primarily of hydrogen, methane, ethane, and propane with various small amounts of nitrogen, hydrogen sulfide, carbon monoxide and hydrocarbons having carbon numbers predominantly in the range of C_4 through C_5.]	272-775-5	68911-58-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-156-00-7	Gases (petroleum), hydrotreated sour kerosine flash drum; Refinery gas; [A complex combination obtained from the flash drum of the unit treating sour kerosine with hydrogen in the presence of a catalyst. It consists primarily of hydrogen and methane with various small amounts of nitrogen, carbon monoxide, and hydro-carbons having carbon numbers predominantly in the range of C_2 through C_5.]	272-776-0	68911-59-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-157-00-2	Gases (petroleum), distillate unifiner desulfurization stripper off; Refinery gas; [A complex combination stripped from the liquid product of the unifiner desulfurization process. It consists of hydrogen sulfide, methane, ethane, and propane.]	272-873-8	68919-01-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-158-00-8	Gases (petroleum), fluidized catalytic cracker fractionation off; Refinery gas; [A complex combination produced by the fractionation of the overhead product of the fluidized catalytic cracking process. It consists of hydrogen, hydrogen sulfide, nitrogen, and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-874-3	68919-02-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-159-00-3	Gases (petroleum), fluidized catalytic cracker scrubbing secondary absorber off; Refinery gas; [A complex combination produced by scrubbing the overhead gas from the fluidized catalytic cracker. It consists of hydrogen, nitrogen, methane, ethane and propane.]	272-875-9	68919-03-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-160-00-9	Gases (petroleum), heavy distillate hydrotreater desulfurization stripper off; Refinery gas; [A complex combination stripped from the liquid product of the heavy distillate hydrotreater desulfurization process. It consists of hydrogen, hydrogen sulfide, and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-876-4	68919-04-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-161-00-4	Gases (petroleum), platformer stabilizer off, light ends fractionation; Refinery gas; [A complex combination obtained by the fractionation of the light ends of the platinum reactors of the plattformer unit. It consists of hydrogen, methane, ethane and propane.]	272-880-6	68919-07-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-162-00-X	Gases (petroleum), preflash tower off, crude distn.; Refinery gas; [A complex combination produced from the first tower used in the distillation of crude oil. It consists of nitrogen and saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-881-1	68919-08-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-163-00-5	Gases (petroleum), tar stripper off; Refinery gas; [A complex combination obtained by the fractionation of reduced crude oil. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	272-884-8	68919-11-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-164-00-0	Gases (petroleum), unifiner stripper off; Refinery gas; [A combination of hydrogen and methane obtained by fractionation of the products from the unifiner unit.]	272-885-3	68919-12-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-165-00-6	Tail gas (petroleum), catalytic hydrodesulfurized naphtha separator; Refinery gas; [A complex combination of hydrocarbons obtained from the hydrodesulfurization of naphtha. It consists of hydrogen, methane, ethane, and propane.]	273-173-5	68952-79-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-166-00-1	Tail gas (petroleum), straight-run naphtha hydrodesulfurizer; Refinery gas; [A complex combination obtained from the hydrodesulfurization of straight-run naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	273-174-0	68952-80-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-167-00-7	Gases (petroleum), sponge absorber off, fluidized catalytic cracker and gas oil desulfurizer overhead fractionation; Refinery gas; [A complex combination obtained by the fractionation of products from the fluidized catalytic cracker and gas oil desulfurizer. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	273-269-7	68955-33-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-168-00-2	Gases (petroleum), crude distn. and catalytic cracking; Refinery gas; [A complex combination produced by crude distillation and catalytic cracking processes. It consists of hydrogen, hydrogen sulfide, nitrogen, carbon monoxide and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	273-563-5	68989-88-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-169-00-8	Gases (petroleum), gas oil diethanolamine scrubber off; Refinery gas; [A complex combination produced by desulfurization of gas oils with diethanolamine. It consists predominantly of hydrogen sulfide, hydrogen and aliphatic hydrocarbons having carbon numbers in the range of C_1 through C_5.]	295-397-2	92045-15-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-170-00-3	Gases (petroleum), gas oil hydrodesulfurization effluent; Refinery gas; [A complex combination obtained by separation of the liquid phase from the effluent from the hydrogenation reaction. It consists predominantly of hydrogen, hydrogen sulfide and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_3.]	295-398-8	92045-16-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-171-00-9	Gases (petroleum), gas oil hydrodesulfurization purge; Refinery gas; [A complex combination of gases obtained from the reformer and from the purges from the hydrogenation reactor. It consists predominantly of hydrogen and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	295-399-3	92045-17-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-172-00-4	Gases (petroleum), hydrogenator effluent flash drum off; Refinery gas; [A complex combination of gases obtained from flash of the effluents after the hydrogenation reaction. It consists predominantly of hydrogen and aliphatic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	295-400-7	92045-18-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-173-00-X	Gases (petroleum), naphtha steam cracking high-pressure residual; Refinery gas; [A complex combination obtained as a mixture of the non-condensable portions from the product of a naphtha steam cracking process as well as residual gases obtained during the preparation of subsequent products. It consists predominantly of hydrogen and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5 with which natural gas may also be mixed.]	295-401-2	92045-19-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-174-00-5	Gases (petroleum), residue visbaking off; Refinery gas; A complex combination obtained from viscosity reduction of residues in a furnace. It consists predominantly of hydrogen sulfide and paraffinic and olefinic hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	295-402-8	92045-20-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-175-00-0	Foots oil (petroleum), acid-treated; Foots oil; [A complex combination of hydrocarbons obtained by treatment of Foot's oil with sulfuric acid. It consists predominantly of branched-chain hydrocarbons with carbon numbers predominantly in the range of C_{20} through C_{50}.]	300-225-7	93924-31-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-176-00-6	Foots oil (petroleum), clay-treated; Foots oil; [A complex combination of hydrocarbons obtained by treatment of Foot's oil with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists predominantly of branched chain hydrocarbons with carbon numbers predominantly in the range of C_{20} through C_{50}.]	300-226-2	93924-32-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-177-00-1	Gases (petroleum), C_{3-4}; Petroleum gas; [A complex combination of hydrocarbons produced by distillation of products from the cracking of crude oil. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_4, predominantly of propane and propylene, and boiling in the range of approximately - 51 °C to - 1 °C (-60 °F to 30 °F.)]	268-629-5	68131-75-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-178-00-7	Tail gas (petroleum), catalytic cracked distillate and catalytic cracked naphtha fractionation absorber; Petroleum gas; [The complex combination of hydrocarbons from the distillation of the products from catalytic cracked distillates and catalytic cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4.]	269-617-2	68307-98-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-179-00-2	Tail gas (petroleum), catalytic polymn. naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons from the fractionation stabilization products from polymerization of naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_1 through C_4.]	269-618-8	68307-99-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-180-00-8	Tail gas (petroleum), catalytic reformed naphtha fractionation stabilizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation stabilization of catalytic reformed naphtha and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-619-3	68308-00-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-181-00-3	Tail gas (petroleum), cracked distillate hydrotreater stripper; Petroleum gas; [A complex combination of hydrocarbons obtained by treating thermal cracked distillates with hydrogen in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-620-9	68308-01-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-182-00-9	Tail gas (petroleum), straight-run distillate hydrodesulfurizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from catalytic hydrodesulfurization of straight run distillates and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-630-3	68308-10-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-183-00-4	Tail gas (petroleum), gas oil catalytic cracking absorber; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of products from the catalytic cracking of gas oil. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-623-5	68308-03-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-184-00-X	Tail gas (petroleum), gas recovery plant; Petroleum gas; [A complex combination of hydrocarbons from the distillation of products from miscellaneous hydrocarbon streams. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-624-0	68308-04-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-185-00-5	Tail gas (petroleum), gas recovery plant deethanizer; Petroleum gas; [A complex combination of hydrocarbons from the distillation of products from miscellaneous hydrocarbon streams. It consists of hydrocarbon having carbon numbers predominantly in the range of C_1 through C_4.]	269-625-6	68308-05-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-186-00-0	Tail gas (petroleum), hydrodesulfurized distillate and hydrodesulfurized naphtha fractionator, acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of hydrodesulfurized naphtha and distillate hydrocarbon streams and treated to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-626-1	68308-06-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-187-00-6	Tail gas (petroleum), hydrodesulfurized vacuum gas oil stripper, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from stripping stabilization of catalytic hydrodesulfurized vacuum gas oil and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-627-7	68308-07-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-188-00-1	Tail gas (petroleum), light straight-run naphtha stabilizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation stabilization of light straight run naphtha and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	269-629-8	68308-09-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-189-00-7	Tail gas (petroleum), propane-propylene alkylation feed prep deethanizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of the reaction products of propane with propylene. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-631-9	68308-11-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-190-00-2	Tail gas (petroleum), vacuum gas oil hydrodesulfurizer, hydrogen sulfide-free; Petroleum gas; [A complex combination of hydrocarbons obtained from catalytic hydrodesulfurization of vacuum gas oil and from which hydrogen sulfide has been removed by amine treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_6.]	269-632-4	68308-12-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-191-00-8	Gases (petroleum), catalytic cracked overheads; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from the catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5 and boiling in the range of approximately - 48 °C to 32 °C (-54 °F to 90 °F).]	270-071-2	68409-99-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-193-00-9	Alkanes, C_{1-2}; Petroleum gas	270-651-5	68475-57-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-194-00-4	Alkanes, C_{2-3}; Petroleum gas	270-652-0	68475-58-1	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-195-00-X	Alkanes, C_{3-4}; Petroleum gas	270-653-6	68475-59-2	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-196-00-5	Alkanes, C_{4-5}; Petroleum gas	270-654-1	68475-60-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-197-00-0	Fuel gases; Petroleum gas; [A combination of light gases. It consists predominantly of hydrogen and/or low molecular weight hydrocarbons.]	270-667-2	68476-26-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-198-00-6	Fuel gases, crude oil of distillates; Petroleum gas; [A complex combination of light gases produced by distillation of crude oil and by catalytic reforming of naphtha. It consists of hydrogen and hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4 and boiling in the range of approximately - 217 °C to - 12 °C (- 423 °F to 10 °F).]	270-670-9	68476-29-9	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-199-00-1	Hydrocarbons, C_{3-4}; Petroleum gas	270-681-9	68476-40-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-200-00-5	Hydrocarbons, C_{4-5}; Petroleum gas	270-682-4	68476-42-6	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-201-00-0	Hydrocarbons, C_{2-4}, C_3-rich; Petroleum gas	270-689-2	68476-49-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-202-00-6	Petroleum gases, liquefied; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_7 and boiling in the range of approximately - 40 °C to 80 °C (- 40 °F to 176 °F).]	270-704-2	68476-85-7	F+; R12 Carc. Cat. 1; R45 Muta. Cat. 2; R46	F+; T R: 12-45-46 S: 53-45		HKS
649-203-00-1	Petroleum gases, liquefied, sweetened; Petroleum gas; [A complex combination of hydrocarbons obtained by subjecting liquefied petroleum gas mix to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_7 and boiling in the range of approximately - 40 °C to 80 °C (-40 °F to 176 °F).]	270-705-8	68476-86-8	F+; R12 Carc. Cat. 1; R45 Muta. Cat. 2; R46	F+; T R: 12-45-46 S: 45-53		HKS
649-204-00-7	Gases (petroleum), C_{3-4}, isobutane-rich; Petroleum gas; [A complex combination of hydrocarbons from the distillation of saturated and unsaturated hydrocarbons usually ranging in carbon numbers from C_3 through C_6, predominantly butane and isobutane. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_4, predominantly isobutane.]	270-724-1	68477-33-8	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-205-00-2	Distillates (petroleum), C_{3-6}, piperylene-rich; Petroleum gas; [A complex combination of hydrocarbons from the distillation of saturated and unsaturated aliphatic hydrocarbons usually ranging in the carbon numbers C_3 through C_6. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly piperylenes.]	270-726-2	68477-35-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-206-00-8	Gases (petroleum), butane splitter overheads; Petroleum gas; [A complex combination of hydrocarbons obtained from the distillation of the butane stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_4.]	270-750-3	68477-69-0	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-207-00-3	Gases (petroleum), C_2-; Petroleum gas; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic fractionation process. It contains predominantly ethane, ethylene, propane, and propylene.]	270-751-9	68477-70-3	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-208-00-9	Gases (petroleum), catalytic-cracked gas oil depropanizer bottoms, C_4-rich acid-free; Petroleum gas; [A complex combination of hydrocarbons obtained from fractionation of catalytic cracked gas oil hydrocarbon stream and treated to remove hydrogen sulfide and other acidic components. It consists of hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly C_4.]	270-752-4	68477-71-4	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-209-00-4	Gases (petroleum), catalytic-cracked naphtha debutanizer bottoms, C_{3-5}-rich; Petroleum gas; [A complex combination of hydrocarbons obtained from the stabilization of catalytic cracked naphtha. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_5.]	270-754-5	68477-72-5	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K
649-210-00-X	Tail gas (petroleum), isomerized naphtha fractionation stabilizer; Petroleum gas; [A complex combination of hydrocarbons obtained from the fractionation stabilization products from isomerized naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_4.]	269-628-2	68308-08-7	⊗ Carc. Cat. 1; R45 Muta. Cat. 2; R46	T R: 45-46 S: 53-45		H K

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649-211-00-5	Foots oil (petroleum), carbon-treated; Foots oil; [A complex combination of hydrocarbons obtained by the treatment of Foots oil with activated carbon for the removal of trace constituents and impurities. It consists predominantly of saturated straight chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-126-0	97862-76-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-212-00-0	Distillates (petroleum), sweetened middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-088-7	64741-86-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-213-00-6	Gas oils (petroleum), solvent-refined; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-092-9	64741-90-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-214-00-1	Distillates (petroleum), solvent-refined middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-093-4	64741-91-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-215-00-7	Gas oils (petroleum), acid-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-112-6	64742-12-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

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649-216-00-2	Distillates (petroleum), acid-treated middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{20} and boiling in the range of approximately 205 °C to 345 °C (401 °F to 653 °F).]	265-113-1	64742-13-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-217-00-8	Distillates (petroleum), acid-treated light; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-114-7	64742-14-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-218-00-3	Gas oils (petroleum), chemically neutralized; Gasoil — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-129-9	64742-29-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-219-00-9	Distillates (petroleum), chemically neutralized middle; Gasoil — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{20} and boiling in the range of approximately 205 °C to 345 °C (401 °F to 653 °F).]	265-130-4	64742-30-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-220-00-4	Distillates (petroleum), clay-treated middle; Gasoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 150 °C to 345 °C (302 °F to 653 °F).]	265-139-3	64742-38-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

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649-221-00-X	Distillates (petroleum), hydrotreated middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-148-2	64742-46-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-222-00-5	Gas oils (petroleum), hydrodesulfurized; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{25} and boiling in the range of approximately 230 °C to 400 °C (446 °F to 752 °F).]	265-182-8	64742-79-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-223-00-0	Distillates (petroleum), hydrodesulfurized middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	265-183-3	64742-80-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-224-00-6	Fuels, diesel; Gasoil — unspecified; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{20} and boiling in the range of approximately 163 °C to 357 °C (325 °F to 675 °F).]	269-822-7	68334-30-5	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H N
649-225-00-1	Fuel oil, No 2; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 32,6 SUS at 37,7 °C (100 °F) to a maximum of 37,9 SUS at 37,7 °C (100 °F).]	270-671-4	68476-30-2	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
649-226-00-7	Fuel oil, No 4; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 45 SUS at 37,7 °C (100 °F) to a maximum of 125 SUS at 37,7 °C (100 °F).]	270-673-5	68476-31-3	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-227-00-2	Fuels, diesel, No 2; Gasoil — unspecified; [A distillate oil having a minimum viscosity of 32,6 SUS at 37,7 °C (100 °F).]	270-676-1	68476-34-6	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
649-228-00-8	Distillates (petroleum), catalytic reformer fractionator residue, high-boiling; Gasoil — unspecified; [A complex combination of hydrocarbons from the distillation of catalytic reformer fracftionator residue. It boils in the range of approximately 343 °C to 399 °C (650 °F to 750 °F).]	270-719-4	68477-29-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-229-00-3	Distillates (petroleum), catalytic reformer fractionator residue, intermediate-boiling; Gasoil — unspecified; [A complex combination of hydrocarbons from the distillation of catalytic reformer fractionator residue. It boils in the range of approximately 288 °C to 371 °C (550 °F to 700 °F).]	270-721-5	68477-30-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-230-00-9	Distillates (petroleum), catalytic reformer fractionator residue, low-boiling; Gasoil — unspecified; [The complex combination of hydrocarbons from the distillation of catalytic reformer fractionator residue. It boils approximately below 288 °C (550 °F).]	270-722-0	68477-31-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-231-00-4	Distillates (petroleum), highly refined middle; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the subjection of a petroleum fraction to several of the following steps: filtration, centrifugation, atmospheric distillation, vacuum distillation, acidification, neutralization and clay treatment. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{20}.]	292-615-8	90640-93-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-232-00-X	Distillates (petroleum) catalytic reformer, heavy arom. conc.; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from the distillation of a catalytically reformed petroleum cut. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{16} and boiling in the range of approximately 200 °C to 300 °C (392 °F to 572 °F).]	295-294-2	91995-34-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-233-00-5	Gas oils, paraffinic; Gasoil — unspecified; [A distillate obtained from the redistillation of a complex combination of hydrocarbons obtained by the distillation of the effluents from a severe catalytic hydrotreatment of paraffins. It boils in the range of approximately 190 °C to 330 °C (374 °F to 594 °F).]	300-227-8	93924-33-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-234-00-0	Naphtha (petroleum), solvent-refined hydrodesulfurized heavy; Gasoil — unspecified	307-035-3	97488-96-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-235-00-6	Hydrocarbons, C_{16-20}, hydrotreated middle distillate, distn. lights; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of a middle distillate with hydrogen. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{20} and boiling in the range of approximately 290 °C to 350 °C (554 °F to 662 °F). It produces a finished oil having a viscosity of 2cSt at 100 °C (212 °F).]	307-659-6	97675-85-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-236-00-1	Hydrocarbons, C_{12-20}, hydrotreated paraffinic, distn. lights; Gasoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of heavy paraffins with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{20} and boiling in the range of approximately 230 °C to 350 °C (446 °F to 662 °F). It produces a finished oil having a viscosity of 2cSt at 100 °C (212 °F).]	307-660-1	97675-86-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-237-00-7	Hydrocarbons, C_{11-17}, solvent-extd. light naphthenic; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a visciosity of 2.2 cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{17} and boiling in the range of approximately 200 °C to 300 °C (392 °F to 572 °F).]	307-757-9	97722-08-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-238-00-2	Gas oils, hydrotreated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained from the redistillation of the effluents from the treatment of paraffins with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{27} and boiling in the range of approximately 330 °C to 340 °C (626 °F to 644 °F).]	308-128-1	97862-78-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-239-00-8	Distillates (petroleum), carbon-treated light paraffinic; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of a petroleum oil fraction with activated charcoal for the removal of traces of polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{28}.]	309-667-5	100683-97-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-240-00-3	Distillates (petroleum), intermediate paraffinic, carbon-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of petroleum with activated charcoal for the removal of trace polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	309-668-0	100683-98-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-241-00-9	Distillates (petroleum), intermediate paraffinic, clay-treated; Gasoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of petroleum with bleaching earth for the removal of trace polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	309-669-6	100683-99-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-242-00-4	Alkanes, C_{12-26}-branched and linear	292-454-3	90622-53-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-243-00-X	Lubricating greases; Grease; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{50}. May contain organic salts of alkali metals, alkaline earth metals, and/or aluminium compounds.]	278-011-7	74869-21-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-244-00-5	Slack wax (petroleum); Slack wax; [A complex combination of hydrocarbons obtained from a petroleum fraction by solvent crystallization (solvent dewaxing) or as a distillation fraction from a very waxy crude. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	265-165-5	64742-61-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-245-00-0	Slack wax (petroleum), acid-treated; Slack wax; [A complex combination of hydrocarbons obtained as a raffinate by treatment of a petroleum slack wax fraction with sulfuric acid treating process. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	292-659-8	90669-77-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-246-00-6	Slack wax (petroleum), clay-treated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of a petroleum slack wax fraction with natural or modified clay in either a contacting or percolation process. It consists predominantly of saturated straight and branched hydrocarbons having carbon numbers predominantly greater than C_{20}.]	292-660-3	90669-78-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-247-00-1	Slack wax (petroleum), hydrotreated; Slack wax; [A complex combination of hydrocarbons obtained by treating slack wax with hydrogen in the presence of a catalyst. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{20}.]	295-523-6	92062-09-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-248-00-7	Slack wax (petroleum), low-melting; Slack wax; [A complex combination of hydrocarbons obtained from a petroleum fraction by solvent deparaffination. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	295-524-1	92062-10-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-249-00-2	Slack wax (petroleum), low-melting, hydrotreated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of low-melting petroleum slack wax with hydrogen in the presence of a catalyst. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C12.]	295-525-7	92062-11-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-250-00-8	Slack wax (petroleum), low-melting, carbon-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting slack wax with activated carbon for the removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-155-9	97863-04-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-251-00-3	Slack wax (petroleum), low-melting, clay-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting petroleum slack wax with bentonite for removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-156-4	97863-05-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-252-00-9	Slack wax (petroleum), low-melting, silicic acid-treated; Slack wax; [A complex combination of hydrocarbons obtained by the treatment of low-melting petroleum slack wax with silicic acid for the removal of trace polar constituents and impurities. It consists predominantly of saturated straight and branched chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-158-5	97863-06-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-253-00-4	Slack wax (petroleum), carbon-treated; Slack wax; [A complex combination of hydrocarbons obtained by treatment of petroleum slack wax with activated charcoal for the removal of trace polar constituents and impurities.]	309-723-9	100684-49-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-254-00-X	Petrolatum; Petrolatum; [A complex combination of hydrocarbons obtained as a semi-solid from dewaxing paraffinic residual oil. It consists predominantly of saturated crystalline and liquid hydrocarbons having carbon numbers predominantly greater than C_{25}.]	232-373-2	8009-03-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-255-00-5	Petrolatum (petroleum), oxidized; Petrolatum; [A complex combination of organic compounds, predominantly high molecular weight carboxylic acids, obtained by the air oxidation of petrolatum.]	265-206-7	64743-01-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-256-00-0	Petrolatum (petroleum), alumina-treated; Petrolatum; [A complex combination of hydrocarbons obtained when petrolatum is treated with Al_2O_3 to remove polar components and impurities. It consists predominantly of saturated, crystalline, and liquid hydrocarbons having carbon numbers predominantly greater than C_{25}.]	285-098-5	85029-74-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-257-00-6	Petrolatum (petroleum), hydrotreated; Petrolatum; [A complex combination of hydrocarbons obtained as a semi-solid from dewaxed paraffinic residual oil treated with hydrogen in the presence of a catalyst. It consists predominantly of saturated microcrystalline and liquid hydrocarbons having carbon numbers predominantly greater than C_{20}.]	295-459-9	92045-77-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-258-00-1	Petrolatum (petroleum), carbon-treated; Petrolatum; [A complex combination of hydrocarbons obtained by the treatment of petroleum petrolatum with activated carbon for the removal of trace polar constituents and impurities. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly greater than C_{20}.]	308-149-6	97862-97-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-259-00-7	Petrolatum (petroleum), silicic acid-treated; Petrolatum; [A complex combination of hydrocarbons obtained by the treatment of petroleum petrolatum with silicic acid for the removal of trace polar constituents and impurities. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly greater than C_{20}.]	308-150-1	97862-98-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-260-00-2	Petrolatum (petroleum), clay-treated; Petrolatum; [A complex combination of hydrocarbons obtained by treatment of petrolatum with bleaching earth for the removal of traces of polar constituents and impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of greater than C_{25}.]	309-706-6	100684-33-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H N
649-261-00-8	Gasoline, natural; Low boiling point naphtha; [A complex combination of hydrocarbons separated from natural gas by processes such as refrigeration or absorption. It consists predominantly of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_8 and boiling in the range of approximately minus 20 °C to 120 °C (-4 °F to 248 °F).]	232-349-1	8006-61-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-262-00-3	Naphtha; Low boiling point naphtha; [Refined, partly refined, or unrefined petroleum products by the distillation of natural gas. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_6 and boiling in the range of approximately 100 °C to 200 °C (212 °F to 392 °F).]	232-443-2	8030-30-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-263-00-9	Ligroine; Low boiling point naphtha; [A complex combination of hydrocarbons obtained by the fractional distillation of petroleum. This fraction boils in a range of approximately 20 °C to 135 °C (58 °F to 275 °F).]	232-453-7	8032-32-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-264-00-4	Naphtha (petroleum), heavy straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-041-0	64741-41-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-265-00-X	Naphtha (petroleum), full-range straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 220 °C (-4 °F to 428 °F).]	265-042-6	64741-42-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-266-00-5	Naphtha (petroleum), light straight-run; Low boiling point naphtha; [A complex combination of hydrocarbons produced by distillation of crude oil. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10} and boiling in the range of approximately minus 20 °C to 180 °C (-4 °F to 356 °F).]	265-046-8	64741-46-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-267-00-0	Solvent naphtha (petroleum), light aliph.; Low boiling point naphtha; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{10} and boiling in the range of approximately 35 °C to 160 °C (95 °F to 320 °F).]	265-192-2	64742-89-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-268-00-6	Distillates (petroleum), straight-run light; Low boiling point naphtha; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_7 and boiling in the range of approximately - 88 °C to 99 °C (-127 °F to 210 °F).]	270-077-5	68410-05-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-269-00-1	Gasoline, vapor-recovery; Low boiling point naphtha; [A complex combination of hydrocarbons separated from the gases from vapor recovery systems by cooling. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately - 20 °C to 196 °C (-4 °F to 384 °F).]	271-025-4	68514-15-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-270-00-7	Gasoline, straight-run, topping-plant; Low boiling point naphtha; [A complex combination of hydrocarbons produced from the topping plant by the distillation of crude oil. It boils in the range of approximately 36,1 °C to 193,3 °C (97 °F to 380 °F).]	271-727-0	68606-11-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-271-00-2	Naphtha (petroleum), unsweetened; Low boiling point naphtha; [A complex combination of hydrocarbons produced from the distillation of naphtha streams from various refinery processes. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 0 °C to 230 °C (25 °F to 446 °F).]	272-186-3	68783-12-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-272-00-8	Distillates (petroleum), light straight-run gasoline fractionation stabilizer overheads; Low boiling point naphtha; [A complex combination of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6.]	272-931-2	68921-08-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-273-00-3	Naphtha (petroleum), heavy straight run, arom.-contg.; Low boiling point naphtha; [A complex combination of hydrocarbons obtained from a distillation process of crude petroleum. It consists predominantly of hydrocarbons having carbon numbers in the range of C_8 through C_{12} and boiling in the range of approximately 130 °C to 210 °C (266 °F to 410 °F).]	309-945-6	101631-20-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-274-00-9	Naphtha (petroleum), full-range alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consist of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 220 °C (194 °F to 428 °F).]	265-066-7	64741-64-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-275-00-4	Naphtha (petroleum), heavy alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 to C_5. It consists of predominantly branched chain saturated hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{12} and boiling in the range of approximately 150 °C to 220 °C (302 °F to 428 °F).]	265-067-2	64741-65-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-276-00-X	Naphtha (petroleum), light alkylate; Low boiling point modified naphtha; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_7 through C_{10} and boiling in the range of aproximately 90 °C to 160 °C (194 °F to 320 °F).]	265-068-8	64741-66-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-277-00-5	Naphtha (petroleum), isomerization; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained from catalytic isomerization of straight chain paraffinic C_4 through C_6 hydrocarbons. It consists predominantly of saturated hydrocarbons such as isobutane, isopentane, 2,2-dimethyl-butane, 2-methylpentane, and 3-methylpentane.]	265-073-5	64741-70-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-278-00-0	Naphtha (petroleum), solvent-refined light; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F).]	265-086-6	64741-84-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-279-00-6	Naphtha (petroleum), solvent-refined heavy; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-095-5	64741-92-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-280-00-1	Raffinates (petroleum), catalytic reformer ethylene glycol-water countercurrent exts.; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinate from the UDEX extraction process on the catalytic reformer stream. It consists of saturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_9.]	270-088-5	68410-71-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-281-00-7	Raffinates (petroleum), reformer, Lurgi unit-sepd.; Low boiling point modified naphtha; [The complex combination of hydrocarbons obtained as a raffinate from a Lurgi separation unit. It consists predominantly of non-aromatic hydrocarbons with various small amounts of aromatic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8.]	270-349-3	68425-35-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-282-00-2	Naphtha (petroleum), full-range alkylate, butane-contg.; Low boiling point modified naphta; [A complex combination of hydrocarbons produced by the distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} with some butanes and boiling in the range of approximately 35 °C to 200 °C (95 °F to 428 °F).]	271-267-0	68527-27-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-283-00-8	Distillates (petroleum), naphtha steam cracking-derived, solvent-refined light hydrotreated; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained as the raffinates from a solvent extraction process of hydrotreated light distillate from steam-cracked naphtha.]	295-315-5	91995-53-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-284-00-3	Naphtha (petroleum), C_{4-12} butane-alkylate, isooctane-rich; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by alkylation of butanes. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12}, rich in isooctane, and boiling in the range of approximately 35 °C to 210 °C (95 °F to 410 °F).]	295-430-0	92045-49-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-285-00-9	Hydrocarbons, hydrotreated light naphtha distillates, solvent-refined; Low boiling point modified naphtha; [A combination of hydrocarbons obtained from the distillation of hydrotreated naphtha followed by a solvent extraction and distillation process. It consists predominantly of saturated hydrocarbons boiling in the range of approximately 94 °C to 99 °C (201 °F to 210 °F).]	295-436-3	92045-55-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-286-00-4	Naphtha (petroleum), isomerization, C_6-fraction; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by distillation of a gasoline which has been catalytically isomerized. It consists predominantly of hexane isomers boiling in the range of approximately 60 °C to 66 °C (140 °F to 151 °F).]	295-440-5	92045-58-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-287-00-X	Hydrocarbons, C_{6-7}, naphtha-cracking, solvent-refined; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by the sorption of benzene from a catalytically fully hydrogenated benzene-rich hydrocarbon cut that was distillatively obtained from prehydrogenated cracked naphtha. It consists predominantly of paraffinic and naphthenic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_7 and boiling in the range of approximately 70 °C to 100 °C (158 °F to 212 °F).]	295-446-8	92045-64-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-288-00-5	Hydrocarbons, C_6-rich, hydrotreated light naphtha distillates, solvent-refined; Low boiling point modified naphtha; [A complex combination of hydrocarbons obtained by distillation of hydrotreated naphtha followed by solvent extraction. It consists predominantly of saturated hydrocarbons and boiling in the range of approximately 65 °C to 70 °C (149 °F to 158 °F).]	309-871-4	101316-67-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-289-00-0	Naphtha (petroleum), heavy catalytic cracked; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by a distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (148 °F to 446 °F). It contains a relatively large proportion of unsaturated hydrocarbons.]	265-055-7	64741-54-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-290-00-6	Naphtha (petroleum), light catalytic cracked; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (-4 °F to 374 °F). It contains a relatively large proportion of unsaturated hydrocarbons.]	265-056-2	64741-55-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-291-00-1	Hydrocarbons, C_{3-11}, catalytic cracker distillates; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillations of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_{11} and boiling in a range approximately up to 204 °C (400 °F).]	270-686-6	68476-46-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-292-00-7	Naphtha (petroleum), catalytic cracked light distd.; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_1 through C_5.]	272-185-8	68783-09-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-293-00-2	Distillates (petroleum), naphtha steam cracking-derived, hydrotreated light arom.; Low boiling point cat-cracked naphtha.; [A complex combination of hydrocarbons obtained by treating a light distillate from steam-cracked naphtha. It consists predom-inantly of aromatic hydrocarbons.]	295-311-3	91995-50-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-294-00-8	Naphtha (petroleum), heavy catalytic cracked, sweetened; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting a catalytic cracked petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 60 °C to 200 °C (140 °F to 392 °F).]	295-431-6	92045-50-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-295-00-3	Naphtha (petroleum), light catalytic cracked sweetened; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting naphtha from a catalytic cracking process to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons boiling in a range of approxim-ately 35 °C to 210 °C (95 °F to 410 °F).]	295-441-0	92045-59-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-296-00-9	Hydrocarbons, C_{8-12}, catalytic-cracking, chem. neutralized; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of a cut from the catalytic cracking process, having undergone an alkaline washing. It consists predominantly of hydrocarbons having carbon numbers in the range of C_8 through C_{12} and boiling in the range of approximately 130 °C to 210 °C (266 °F to 410 °F).]	295-794-0	92128-94-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-297-00-4	Hydrocarbons, C_{8-12}, catalytic cracker distillates; Low boiling point cat-cracked naphtha; [A complex combination of hydrocarbons obtained by distillation of products from a catalytic cracking process. It consists pre-dominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12} and boiling in the range of approximately 140 °C to 210 °C (284 °F to 410 °F).]	309-974-4	101794-97-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-298-00-X	Hydrocarbons, C_{8-12}, catalytic cracking, chem. neutralized, sweetened; Low boiling point cat-cracked naphtha	309-987-5	101896-28-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-299-00-5	Naphtha (petroleum), light catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced from the distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F). It contains a relatively large proportion of aromatic and branched chain hydrocarbons. This stream may contain 10 vol. % or more benzene.]	265-065-1	64741-63-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-300-00-9	Naphtha (petroleum), heavy catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced from the distillation of products from a catalytic reforming process. It consists of predominantly aromatic hydrocarbons having numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-070-9	64741-68-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-301-00-4	Distillates (petroleum), catalytic reformed depentanizer; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons from the distillation of products from a catalytic reforming process. It consists predominantly of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 49 °C to 63 °C - 57 °F to 145 °F).]	270-660-4	68475-79-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-302-00-X	Hydrocarbons, C_{2-6}, C_{6-8} catalytic reformer; Low boiling point cat-reformed naphtha	270-687-1	68476-47-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-303-00-5	Residues (petroleum), C_{6-8} catalytic reformer; Low boiling point cat-reformed naphtha; [A complex residuum from the catalytic reforming of C_{6-8} feed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	270-794-3	68478-15-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-304-00-0	Naphtha (petroleum), light catalytic reformed, arom.-free; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained from distillation of products from a catalytic reforming process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_8 and boiling in the range of approximately 35 °C to 120 °C (95 °F to 248 °F). It contains a relatively large proportion of branched chain hydrocarbons with the aromatic components removed.]	270-993-5	68513-03-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-305-00-6	Distillates (petroleum), catalytic reformed straight-run naphtha overheads; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by the catalytic reforming of straight-run naphtha followed by the fractionation of the total effluent. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	271-008-1	68513-63-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-306-00-1	Petroleum products, hydrofiner-powerformer reformates; Low boiling point cat-reformed naphtha; [The complex combination of hydrocarbons obtained in a hydrofiner-powerformer process and boiling in a range of approximately 27 °C to 210 °C (80 °F to 410 °F).]	271-058-4	68514-79-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-307-00-7	Naphtha (petroleum, full-range reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced by the distillation of the products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 35 °C to 230 °C (95 °F to 446 °F).]	272-895-8	68919-37-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-308-00-2	Naphtha (petroleum), catalytic reformed; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic reforming process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 30 °C to 220 °C (90 °F to 430 °F). It contains a relatively large proportion of aromatic and branched chain hydrocarbons. This stream may contain 10 vol. % or more benzene.]	273-271-8	68955-35-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-309-00-8	Distillates (petroleum), catalytic reformed hydrotreated light, C_{8-12} arom. fraction; Low boiling point cat-reformed naphtha; [A complex combination of alkylbenzenes obtained by the catalytic reforming of petroleum naphtha. It consists predominantly of alkylbenzenes having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 160 °C to 180 °C (320 °F to 356 °F).]	285-509-8	85116-58-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-310-00-3	Aromatic hydrocarbons, C_8, catalytic reforming-derived; Low boiling point cat-reformed naphtha	295-279-0	91995-18-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-311-00-9	Aromatic hydrocarbons, C_{7-12}, C_8-rich; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} (primarily C_8) and can contain nonaromatic hydrocarbons, both boiling in the range of approximately 130 °C to 200 °C (266 °F to 392 °F).]	297-401-8	93571-75-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-312-00-4	Gasoline, C_{5-11}, high-octane stabilized reformed; Low boiling point cat-reformed naphtha; [A complex high octane combination of hydrocarbons obtained by the catalytic dehydrogenation of a predominantly naphthenic naphtha. It consists predominantly of aromatics and non-aromatics having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 45 °C to 185 °C (113 °F to 365 °F).]	297-458-9	93572-29-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-313-00-X	Hydrocarbons, C_{7-12}, C $_{>9}$-arom.-rich, reforming heavy fraction; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of nonaromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 120 °C to 210 °C (248 °F to 380 °F) and C_9 and higher aromatic hydrocarbons.]	297-465-7	93572-35-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-314-00-5	Hydrocarbons, C_{5-11}, nonaroms.-rich, reforming light fraction; Low boiling point cat-reformed naphtha; [A complex combination of hydrocarbons obtained by separation from the platformate-containing fraction. It consists predominantly of nonaromatic hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 35 °C to 125 °C (94 °F to 257 °F), benzene and toluene.]	297-466-2	93572-36-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-315-00-0	Foots oil (petroleum), silicic acid-treated; Foots oil; [A complex combination of hydrocarbons obtained by the treatment of Foots oil with silicic acid for removal of trace constituents and impurities. It consists predominantly of straight chain hydrocarbons having carbon numbers predominantly greater than C_{12}.]	308-127-6	97862-77-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-316-00-6	Naphtha (petroleum), light thermal cracked; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons from distillation of products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_8 and boiling in the range of approximately minus 10 °C to 130 °C (14 °F to 266 °F).]	265-075-6	64741-74-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-317-00-1	Naphtha (petroleum), heavy thermal cracked; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons from distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 220 °C (148 °F to 428 °F).]	265-085-0	64741-83-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-318-00-7	Distillates (petroleum), heavy arom.; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons from the distillation of the products from the thermal cracking of ethane and propane. This higher boiling fraction consists predominantly of C_5-C_7 aromatic hydrocarbons with some unsaturated aliphatic hydrocarbons having carbon number predominantly of C_5. This stream may contain benzene.]	267-563-4	67891-79-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-319-00-2	Distillates (petroleum), light arom.; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons from the distillation of the products from the thermal cracking of ethane and propane. This lower boiling fraction consists predominantly of C_5-C_7 aromatic hydrocarbons with some unsaturated aliphatic hydrocarbons having a carbon number predominantly of C_5. This stream may contain benzene.]	267-565-5	67891-80-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-320-00-8	Distillates (petroleum), naphtha-raffinate pyrolyzate-derived, gasoline-blending; Low boiling point thermally cracked naphtha; [The complex combination of hydrocarbons obtained by the pyrolysis fractionation at 816 °C (1500 °F) of naphtha and raffinate. It consists predominantly of hydrocarbons having a carbon number of C_9 and boiling at approximately 204 °C (400 °F).]	270-344-6	68425-29-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-321-00-3	Aromatic hydrocarbons, C_{6-8}, naphtha-raffinate pyrolyzate-derived; Low boiling point thermally cracked naphtha; A complex combination of hydrocarbons obtained by the fractionation pyrolysis at 816 °C (1500 °F) of naphtha and raffinate. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8, including benzene.]	270-658-3	68475-70-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-322-00-9	Distillates (petroleum), thermal cracked naphtha and gas oil; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by distillation of thermally cracked naphtha and/or gas oil. It consists predominantly of olefinic hydrocarbons having a carbon number of C_5 and boiling in the range of approximately 33 °C to 60 °C (91 °F to 140 °F).]	271-631-9	68603-00-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-323-00-4	Distillates (petroleum), thermal cracked naphtha and gas oil, C_5-dimer-contg.; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the extractive distillation of thermal cracked naphtha and/or gas oil. It consists predominantly of hydrocarbons having a carbon number of C_5 with some dimerized C_5 olefins and boiling in the range of approximately 33 °C to 184 °C (91 °F to 363 °F).]	271-632-4	68603-01-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-324-00-X	Distillates (petroleum), thermal cracked naphtha and gas oil, extractive; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the extractive distillation of thermal cracked naphtha and/or gas oil. It consists of paraffinic and olefinic hydrocarbons, predominantly isoamylenes such as 2-methyl-1-butene and 2-methyl-2-butene and boiling in the range of approximately 31 °C to 40 °C (88 °F to 104 °F).]	271-634-5	68603-03-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-325-00-5	Distillates (petroleum), light thermal cracked, debutanized arom.; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons produced by the distillation of products from a thermal cracking process. It consists predominantly of aromatic hydrocarbons, primarily benzene.]	273-266-0	68955-29-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-326-00-0	Naphtha (petroleum), light thermal cracked, sweetened; Low boiling point thermally cracked naphtha; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate from the high temperature thermal cracking of heavy oil fractions to a sweetening process to convert mercaptans. It consists predominantly of aromatics, olefins and saturated hydrocarbons boiling in the range of approximately 20 °C to 100 °C (68 °F to 212 °F).]	295-447-3	92045-65-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-327-00-6	Naphtha (petroleum), hydrotreated heavy; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{13} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-150-3	64742-48-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-328-00-1	Naphtha (petroleum), hydrotreated light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-151-9	64742-49-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-329-00-7	Naphtha (petroleum), hydrodesulfurized light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-178-6	64742-73-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-330-00-2	Naphtha (petroleum), hydrodesulfurized heavy; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-185-4	64742-82-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-331-00-8	Distillates (petroleum), hydrotreated middle, intermediate boiling; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by the distillation of products from a middle distillate hydrotreating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{10} and boiling in the range of approximately 127 °C to 188 °C (262 °F to 370 °F).]	270-092-7	68410-96-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-332-00-3	Distillates (petroleum), light distillate hydrotreating process, low-boiling; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by the distillation of products from the light distillate hydrotreating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_9 and boiling in the range of approximately 3 °C to 194 °C (37 °F to 382 °F).]	270-093-2	68410-97-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-333-00-9	Distillates (petroleum), hydrotreated heavy naphtha, deiso-hexanizer overheads; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by distillation of the products from a heavy naphtha hydro-treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 49 °C to 68 °C (- 57 °F to 155 °F).]	270-094-8	68410-98-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-334-00-4	Solvent naphtha (petroleum), light arom., hydrotreated; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F).]	270-988-8	68512-78-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-335-00-X	Naphtha (petroleum), hydrodesulfurized thermal cracked light; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by fractionation of hydrodesulfurized thermal cracker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 23 °C to 195 °C (73 °F to 383 °F).]	285-511-9	85116-60-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-336-00-5	Naphtha (petroleum), hydrotreated light, cycloalkane-contg.; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from the distillation of a petroleum fraction. It consists predominantly of alkanes and cycloalkanes boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	285-512-4	85116-61-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-337-00-0	Naphtha (petroleum), heavy steam-cracked, hydrogenated; Low boiling point hydrogen treated naphtha	295-432-1	92045-51-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-338-00-6	Naphtha (petroleum), hydrodesulfurized full-range; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained from a catalytic hydrodesulfurization process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately 30 °C to 250 °C (86 °F to 482 °F).]	295-433-7	92045-52-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-339-00-1	Naphtha (petroleum), hydrotreated light steam-cracked; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction, derived from a pyrolysis process, with hydrogen in the presence of a catalyst. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{11} and boiling in the range of approximately 35 °C to 190 °C (95 °F to 374 °F).]	295-438-4	92045-57-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-340-00-7	Hydrocarbons, C_{4-12}, naphtha-cracking, hydrotreated; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by distillation from the product of a naphtha steam cracking process and subsequent catalytic selective hydrogenation of gum formers. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 30 °C to 230 °C (86 °F to 446 °F).]	295-443-1	92045-61-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-341-00-2	Solvent naphtha (petroleum), hydrotreated light naphthenic; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists predominantly of cycloparaffinic hydrocarbons having carbon numbers predominantly in the range of C_6 through C_7 and boiling in the range of approximately 73 °C to 85 °C (163 °F to 185 °F).]	295-529-9	92062-15-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-342-00-8	Naphtha (petroleum), light steam-cracked, hydrogenated; Low boiling point hydrogen treated naphtha; [A complex comination of hydrocarbons produced from the separation and subsequent hydrogenation of the products of a steam-cracking process to produce ethylene. It consists predominantly of saturated and unsaturated paraffins, cyclic paraffins and cyclic aromatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10} and boiling in the range of approximately 50 °C to 200 °C (122 °F to 392 °F). The proportion of benzene hydrocarbons may vary up to 30 wt. % and the stream may also contain small amounts of sulphur and oxygenated compounds.]	296-942-7	93165-55-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-343-00-3	Hydrocarbons, C_{6-11}, hydrotreated, dearomatized; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-852-0	93763-33-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-344-00-9	Hydrocarbons, C_{9-12}, hydrotreated, dearomatized; Low boiling point hydrogen treated naphtha; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-853-6	93763-34-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-345-00-4	Stoddard solvent; Low boiling point naphtha — unspecified; [A colourless, refined petroleum distillate that is free from rancid or objectionable odors and that boils in a range of approximately 300 °F to 400 °F.]	232-489-3	8052-41-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-346-00-X	Natural gas condensates (petroleum); Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated as a liquid from natural gas in a surface separator by retrograde condensation. It consists mainly of hydrocarbons having carbon numbers predominantly in the range of C_2 to C_{20}. It is a liquid at atmospheric temperature and pressure.]	265-047-3	64741-47-5	Carc. Cat. 2; R4 Xn; R65	T R: 45-65 S: 53-45		H P
649-347-00-5	Natural gas (petroleum), raw liq. mix; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated as a liquid from natural gas in a gas recycling plant by processes such as refrigeration or absorption. It consists mainly of saturated aliphatic hydrocarbons having carbon numbers in the range of C_2 through C_8.]	265-048-9	64741-48-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-348-00-0	Naphtha (petroleum), light hydrocracked; Low boiling naphtha — unspecified; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{10}, and boiling in the range of approximately minus 20 °C to 180 °C (- 4 °F to 356 °F).]	265-071-4	64741-69-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-349-00-6	Naphtha (petroleum), heavy hydrocracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12}, and boiling in the range of approximately 65 °C to 230 °C (148 °F to 446 °F).]	265-079-8	64741-78-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-350-00-1	Naphtha (petroleum), sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately minus 10 °C to 230 °C (14 °F to 446 °F).]	265-089-2	64741-87-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-351-00-7	Naphtha (petroleum), acid-treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as a raffinate from a sulfuric acid treating process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 230 °C (194 °F to 446 °F).]	265-115-2	64742-15-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-352-00-2	Naphtha (petroleum), chemically neutralized heavy; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{12} and boiling in the range of approximately 65 °C to 230 °C (149 °F to 446 °F).]	265-122-0	64742-22-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-353-00-8	Naphtha (petroleum), chemically neutralized light; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F).]	265-123-6	64742-23-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-354-00-3	Naphtha (petroleum), catalytic dewaxed; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the catalytic dewaxing of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12} and boiling in the range of approximately 35 °C to 230 °C (95 °F to 446 °F).]	265-170-2	64742-66-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-355-00-9	Naphtha (petroleum), light steam-cracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of the products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately minus 20 °C to 190 °C (- 4 °F to 374 °F). This stream is likely to contain 10 vol.% or more benzene.]	265-187-5	64742-83-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-356-00-4	Solvent naphtha (petroleum), light arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from distillation of aromatic streams. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{10} and boiling in the range of approximately 135 °C to 210 °C (275 °F to 410 °F).]	265-199-0	64742-95-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-357-00-X	Aromatic hydrocarbons, C_{6-10}, acid-treated, neutralized; Low boiling point naphtha — unspecified	268-618-5	68131-49-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-358-00-5	Distillates (petroleum), C_{3-5}, 2-methyl-2-butene-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons from the distillation of hydrocarbons usually ranging in carbon numbers from C_3 through C_5, predominantly isopentane and 3-methyl-1-butene. It consists of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_3 through C_5, predominantly 2-methyl-2-butene.]	270-725-7	68477-34-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-359-00-0	Distillates (petroleum), polymd. steam-cracked petroleum distillates, C_{5-12} fraction; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the distillation of polymerized steam-cracked petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12}.]	270-735-1	68477-50-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-360-00-6	Distillates (petroleum), steam-cracked, C_{5-12} fraction; Low boiling point naphtha — unspecified; [A complex combination of organic compounds obtained by the distillation of products from a steam cracking process. It consists of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_5 through C_{12}.]	270-736-7	68477-53-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-361-00-1	Distillates (petroleum), steam-cracked, C_{5-10} fraction, mixed with light steam-cracked petroleum naphtha C_5 fraction; Low boiling point naphtha — unspecified	270-738-8	68477-55-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-362-00-7	Extracts (petroleum), cold-acid, C_{4-6}; Low boiling point naphtha — unspecified; [A complex combination of organic compounds produced by cold acid unit extraction of saturated and unsaturated aliphatic hydrocarbons usually ranging in carbon numbers from C_3 through C_6, predominantly pentanes and amylenes. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers in the range of C_4 through C_6, predominantly C_5.]	270-741-4	68477-61-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-363-00-2	Distillates (petroleum), dependentanizer overheads; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from a catalytic cracked gas stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6.]	270-771-8	68477-89-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-364-00-8	Residues (petroleum), butane splitter bottoms; Low boiling point naphtha — unspecified; [A complex residuum from the distillation of butane stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4through C_6.]	270-791-7	68478-12-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-365-00-3	Residual oils (petroleum), deisobutanizer tower; Low boiling point naphtha — unspecified; [A complex residuum from the atmospheric distillation of the butane-butylene stream. It consists of aliphatic hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6.]	270-795-9	68478-16-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-366-00-9	Naphtha (petroleum), full-range coker; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by the distillation of products from a fluid coker. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{15} and boiling in the range of approximately 43 °C to 250 °C (110 °F to 500 °F).]	270-991-4	68513-02-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-367-00-4	Naphtha (petroleum), steam-cracked middle arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by the distillation of products from a steam-cracking process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 130 °C to 220 °C (266 °F to 428 °F).]	271-138-9	68516-20-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-368-00-X	Naphtha (petroleum), clay-treated full-range straight-run; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons resulting from treatment of full-range straight-run naphtha with natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{11} and boiling in the range of approximately - 20 °C to 220 °C (- 4 °F to 429 °F).]	271-262-3	68527-21-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-369-00-5	Naphtha (petroleum), clay-treated light straight-run; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons resulting from treatment of light straight-run naphtha with a natural or modified clay, usually in a percolation process to remove the trace amounts of polar compounds and impurities, present. It consists of hydro-carbons having carbon numbers predominantly in the range of C_7 through C_{10} and boiling in the range of approximately 93 °C to 180 °C (200 °F to 356 °F.]	271-263-9	68527-22-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-370-00-0	Naphtha (petroleum), light steam-cracked arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by distillation of products from a steam-cracking process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_9 and boiling in the range of approximately 110 °C to 165 °C (230 °F to 329 °F).]	271-264-4	68527-23-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-371-00-6	Naphtha (petroleum), light steam-cracked, debenzenized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by distillation of products from a steam-cracking process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_{12} and boiling in the range of approximately 80 °C to 218 °C (176 °F to 424 °F).]	271-266-5	68527-26-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-372-00-1	Naphtha (petroleum), arom.-contg.; Low boiling point naphtha — unspecified	271-635-0	68603-08-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-373-00-7	Gasoline, pyrolysis, debutanizer bottoms; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the fractionation of depropanizer bottoms. It consists of hydrocarbons having carbon numbers predominantly greater than C_5.]	271-726-5	68606-10-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-374-00-2	Naphtha (petroleum), light, sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of saturated and unsaturated hydrocarbons having carbon numbers predominantly in the range of C_3 through C_6 and boiling in the range of approximately - 20 °C to 100 °C (- 4 °F to 212 °F).]	272-206-0	68783-66-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-375-00-8	Natural gas condensates; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons separated and/or condensed from natural gas during transportation and collected at the wellhead and/or from the production, gathering, transmission, and distribution pipelines in deeps, scrubbers, etc. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_2 through C_8.]	272-896-3	68919-39-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H J
649-376-00-3	Distillates (petroleum), naphtha unifiner stripper; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons produced by stripping the products from the naphtha unifiner. It consists of saturated aliphatic hydrocarbons having carbon numbers predominantly in the range of C_2 through C_6.]	272-932-8	68921-09-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-377-00-9	Naphtha (petroleum), catalytic reformed light, arom.-free fraction; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons remaining after removal of aromatic compounds from catalytic reformed light naphtha in a selective absorption process. It consists predominantly of paraffinic and cyclic compounds having carbon numbers predominantly in the range of C_5 to C_8 and boiling in the range of approximately 66 °C to 121 °C (151 °F to 250 °F).]	285-510-3	85116-59-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-378-00-4	Gasoline; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons consisting primarily of paraffins, cycloparaffins, aromatic and olefinic hydrocarbons having carbon numbers predominantly greater than C_3 and boiling in the range of 30 °C to 260 °C (86 °F to 500 °F).]	289-220-8	86290-81-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-379-00-X	Aromatic hydrocarbons, C_{7-8}, dealkylation products, distn. residues; Low boiling point naphtha — unspecified	292-698-0	90989-42-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-380-00-5	Hydrocarbons, C_{4-6}, depentanizer lights, arom. hydrotreater; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the depentanizer column before hydrotreatment of the aromatic charges. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_6, predominantly pentanes and pentenes, and boiling in the range of approximately 25 °C to 40 °C (77 °F to 104 °F).]	295-298-4	91995-38-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-381-00-0	Distillates (petroleum), heat-soaked steam-cracked naphtha, C_5-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of heat-soaked steam-cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_4 through C_6, predominantly C_5.]	295-302-4	91995-41-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-382-00-6	Extracts (petroleum), catalytic reformed light naphtha solvent; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained as the extract from the solvent extraction of a catalytically reformed petroleum cut. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_8 and boiling in the range of approximately 100 °C to 200 °C (212 °F to 392 °F).]	295-331-2	91995-68-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-383-00-1	Naphtha (petroleum), hydrodesulfurized light, dearomatized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of hydrodesulfurized and dearomatized light petroleum fractions. It consists predominantly of C_7 paraffins and cycloparaffins boiling in a range of approximately 90 °C to 100 °C (194 °F to 212 °F).]	295-434-2	92045-53-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-384-00-7	Naphtha (petroleum), light, C_5-rich, sweetened; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists of hydrocarbons having carbon numbers predominantly in the range of C_4 through C_5, predominantly C_5, and boiling in the range of approximately minus 10 °C to 35 °C (14 °F to 95 °F).]	295-442-6	92045-60-8	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-385-00-2	Hydrocarbons, C_{8-11}, naphtha-cracking, toluene cut; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation from prehydrogenated cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{11} and boiling in the range of approximately 130 °C to 205 °C (266 °F to 401 °F).]	295-444-7	92045-62-0	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-386-00-8	Hydrocarbons, C_{4-11}, naphtha-cracking, arom.-free; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from prehydrogenated cracked naphtha after distillative separation of benzene- and toluene-containing hydrocarbon cuts and a higher boiling fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range C_4 through C_{11} and boiling in the range of approximately 30 °C to 205 °C (86 °F to 401 °F).]	295-445-2	92045-63-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-387-00-3	Naphtha (petroleum), light heat-soaked, steam-cracked; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the fractionation of steam cracked naphtha after recovery from a heat soaking process. It consists predominantly of hydrocarbons having a carbon numbers predominantly in the range of C_4 through C_6 and boiling in the range of approximately 0 °C to 80 °C (32 °F to 176 °F.]	296-028-8	92201-97-3	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

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649-388-00-9	Distillates (petroleum), C_6-rich; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained from the distillation of a petroleum feedstock. It consists predominantly of hydrocarbons having carbon numbers of C_5 through C_7, rich in C_6, and boiling in the range of approximately 60 °C to 70 °C (140 °F to 158 °F).]	296-903-4	93165-19-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-389-00-4	Gasoline, pyrolysis, hydrogenated; Low boiling point naphtha-unspecified; [A distillation fraction from the hydrogenation of pyrolysis gasoline boiling in the range of approximately 20 °C to 200 °C (68 °F to 392 °F).]	302-639-3	94114-03-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-390-00-X	Distillates (petroleum), steam-cracked, C_{8-12} fraction, polymd., distn. lights; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of the polymerized C_8 through C_{12} fraction from steam-cracked petroleum distillates. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12}.]	305-750-5	95009-23-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-391-00-5	Extracts (petroleum) heavy naphtha solvent, clay-treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment of heavy naphthic solvent petroleum extract with bleaching earth. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{18} and boiling in the range of approximately 80 °C to 180 °C (175 °F to 356 °F).]	308-261-5	97926-43-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-392-00-0	Naphtha (petroleum), light steam-cracked, debenzenized, thermally treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment and distillation of debenzenized light steam-cracked petroleum naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 95 °C to 200 °C (203 °F to 392 °F).]	308-713-1	98219-46-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-393-00-6	Naphtha (petroleum), light steam-cracked, thermally treated; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the treatment and distillation of light steam-cracked petroleum naphtha. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_6 and boiling in the range of approximately 35 °C to 80 °C (95 °F to 176 °F).]	308-714-7	98219-47-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-394-00-1	Distillates (petroleum), C_{7-9}, C_8-rich, hydrodesulfurized dearomatized; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of petroleum light fraction, hydrodesulfurized and dearomatized. It consists predominantly of hydrocarbons having carbon numbers in the range of C_7 through C_9, predominantly C_8 paraffins and cycloparaffins, boiling in the range of approximately 120 °C to 130 °C (248 °F to 266 °F).]	309-862-5	101316-56-7	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-395-00-7	Hydrocarbons, C_{6-8}, hydrogenated sorption-dearomatized, toluene raffination; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained during the sorptions of toluene from a hydrocarbon fraction from cracked gasoline treated with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_6 through C_8 and boiling in the range of approximately 80 °C to 135 °C (176 °F to 275 °F).]	309-870-9	101316-66-9	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-396-00-2	Naphtha (petroleum), hydrodesulfurized full-range coker; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized coker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 to C_{11} and boiling in the range of approximately 23 °C to 196 °C (73 °F to 385 °F).]	309-879-8	101316-76-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-397-00-8	Naphtha (petroleum), sweetened light; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum naphtha to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_5 through C_8 and boiling in the range of approximately 20 °C to 130 °C (68 °F to 266 °F).]	309-976-5	101795-01-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-398-00-3	Hydrocarbons, C_{3-6}, C_5-rich, steam-cracked naphtha; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of steam-cracked naphtha. It consists predominantly of hydrocarbons having carbon numbers in the range of C_3 through C_6, predominantly C_5.]	310-012-0	102110-14-5	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-399-00-9	Hydrocarbons, C_5-rich, dicyclopentadiene-contg.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by distillation of the products from a stream-cracking process. It consists predominantly of hydrocarbons having carbon numbers of C_5 and dicyclopentadiene and boiling in the range of approximately 30 °C to 170 °C (86 °F to 338 °F).]	310-013-6	102110-15-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-400-00-2	Residues (petroleum), steam-cracked light, arom.; Low boiling point naphtha — unspecified; [A complex combination of hydrocarbons obtained by the distillation of the products of steam cracking or similar processes after taking off the very light products resulting in a residue starting with hydrocarbons having carbon numbers greater than C_5. It consists predominantly of aromatic hydrocarbons having carbon numbers greater than C_5 and boiling above approximately 40 °C (104 °F).]	310-057-6	102110-55-4	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-401-00-8	Hydrocarbons, C ≥ 5, C_{5-6}-rich; Low boiling point naphtha — unspecified	270-690-8	68476-50-6	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-402-00-3	Hydrocarbons, C_5-rich; Low boiling point naphtha — unspecified	270-695-5	68476-55-1	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-403-00-9	Aromatic hydrocarbons, C_{8-10}; Low boiling point naphtha — unspecified	292-695-4	90989-39-2	Carc. Cat. 2; R45 Xn; R65	T R: 45-65 S: 53-45		H P
649-404-00-4	Kerosine (petroleum); Straight run kerosine; [A complex combination of hydrocarbons produced by the distillation of crude oil. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (320 °F to 554 °F).]	232-366-4	8008-20-6	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-405-00-X	Solvent naphtha (petroleum), medium aliph.; Straight run kerosine; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{12} and boiling in the range of approximately 140 °C to 220 °C (284 °F to 428 °F).]	265-191-7	64742-88-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-406-00-5	Solvent naphtha (petroleum) heavy aliph.; Straight run kerosine; [A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{16} and boiling in the range of approximately 190 °C to 290 °C (374 °F to 554 °F).]	265-200-4	64742-96-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-407-00-0	Kerosine (petroleum), straight-run wide-cut; Straight run kerosine; [A complex combination of hydrocarbons obtained as a wide cut hydrocarbon fuel cut from atmospheric distillation and boiling in the range of approximately 70 °C to 220 °C (158 °F to 428 °F).]	295-418-5	92045-37-9	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-408-00-6	Distillates (petroleum), steam-cracked; Cracked kerosine; [A complex combination of hydrocarbons obtained by the distillation of the products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{16} and boiling in the range of approximately 90 °C to 290 °C (190 °F to 554 °F).]	265-194-3	64742-91-2	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-409-00-1	Distillates (petroleum), cracked stripped steam-cracked petroleum distillates, C_{8-10} fraction; Cracked kerosine; [A complex combination of hydrocarbons obtained by distilling cracked stripped steam-cracked distillates. It consists of hydro-carbons having carbon numbers in the range of C_8 through C_{10} and boiling in the range of approximately 129 °C to 194 °C (264 °F to 382 °F).]	270-728-3	68477-39-4	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-410-00-7	Distillates (petroleum), cracked stripped steam-cracked petroleum distillates, C_{10-12} fraction; Cracked kerosine; [A complex combination of hydrocarbons obtained by distilling cracked stripped steam-cracked distillates. It consists predominantly of aromatic hydrocarbons having carbon numbers in the range of C_{10} through C_{12}.]	270-729-9	68477-40-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-411-00-2	Distillates (petroleum), steam-cracked, C_{8-12} fraction; Cracked kerosine; [A complex combination of organic compounds obtained by the distillation of products from a steam cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{12}.]	270-737-2	68477-54-3	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-412-00-8	Kerosine (petroleum), hydrodesulfurized thermal cracked; Cracked kerosine; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized thermal cracker distillate. It consists predominantly of hydrocarbons predominantly in the range of C_8 to C_{16} and boiling in the range of approximately 120 °C to 283 °C (284 °F to 541 °F).]	285-507-7	85116-55-8	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-413-00-3	Aromtic hydrocarbons, $C_{\geq 10}$, steam-cracking, hydrotreated; Cracked kerosine; [A complex combination of hydrocarbons produced by the distillation of the products from a steam cracking process treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly greater than C_{10} and boiling in the range of approximately 150 °C to 320 °C (302 °F to 608 °F).]	292-621-0	90640-98-5	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-414-00-9	Naphtha (petroleum), steam-cracked, hydrotreated, C_{9-10}-arom.-rich; Cracked kerosine; [A complex combination of hydrocarbons produced by the distillation of the products from a steam cracking process thereafter treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers in the range of C_9 through C_{10} and boiling in the range of approximately 140 °C to 200 °C (284 °F to 392 °F).]	292-637-8	90641-13-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-415-00-4	Distillates (petroleum), thermal-cracked, alkylarom. hydrocarbon-rich; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of thermal-cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F.]	309-866-7	101316-61-4	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-416-00-X	Distillates (petroleum), catalytic cracked heavy tar light; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of catalytic cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F).]	309-938-8	101631-13-4	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-417-00-5	Solvent naphtha (petroleum), hydrocracked heavy arom.; Cracked kerosine; [A complex combination of hydrocarbons obtained by the distillation of hydrocracked petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 235 °C to 290 °C (455 °F to 554 °F).]	309-881-9	101316-80-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-418-00-0	Distillates (petroleum), steam-cracked heavy tar light; Cracked kerosine; [A complex combination of hydrocarbons obtained by distillation of steam cracking heavy tars. It consists predominantly of highly alkylated aromatic hydrocarbons boiling in the range of approximately 100 °C to 250 °C (212 °F to 482 °F).]	309-940-9	101631-15-6	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-419-00-6	Distillates (petroleum), alkylate; Kerosine — unspecified; [A complex combination of hydrocarbons produced by distillation of the reaction products of isobutane with monoolefinic hydrocarbons usually ranging in carbon numbers from C_3 through C_5. It consists of predominantly branched chain saturated hydro-carbons having carbon numbers predominantly in the range of C_{11} through C_{17} and boiling in the range of approximately 205 °C to 320 °C (401 °F to 608 °F).]	265-074-0	64741-73-7	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-420-00-1	Extracts (petroleum), heavy naphtha solvent; Kerosine — unspecified; [A complex combination of hydrocarbons obtained as the extract from a solvent extraction process. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_7 through C_{12} and boiling in the range of approximately 90 °C to 220 °C (194 °F to 428 °F).]	265-099-7	64741-98-6	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-421-00-7	Distillates (petroleum), chemically neutralized light; Kerosine — unspecified; [A complex combination of hydrocarbons produced by a treating process to remove acidic materials. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-132-5	64742-31-0	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-422-00-2	Distillates (petroleum), hydrotreated light; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-149-8	64742-47-8	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-423-00-8	Kerosine (petroleum), hydrodesulfurized; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 150 °C to 290 °C (302 °F to 554 °F).]	265-184-9	64742-81-0	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-424-00-3	Solvent naphtha (petroleum), heavy arom.; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from distillation of aromatic streams. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of approximately 165 °C to 290 °C (330 °F to 554 °F).]	265-198-5	64742-94-5	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-425-00-9	Naphtha (petroleum), heavy coker; Kerosine — unspecified; [A complex combination of hydrocarbons from the distillation of products from a fluid coker. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_6 through C_{15} and boiling in the range of approximately 157 °C to 288 °C (315 °F to 550 °F).]	269-778-9	68333-23-3	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-426-00-4	Naphtha (petroleum), catalytic reformed hydrodesulfurized heavy, arom. fraction; Kerosine — unspecified; [A complex combination of hydrocarbons produced by fractionation from catalytically reformed hydrodesulfurized naphtha. It consists predominantly of aromatic hydrocarbons having carbon numbers predominently in the range of C_7 to C_{13} and boiling in the range of approximately 98 °C to 218 °C (208 °F to 424 °F).]	285-508-2	85116-57-0	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-427-00-X	Kerosine (petroleum), sweetened; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by subjecting a petroleum distillate to a sweetening process to convert mercaptans or to remove acidic impurities. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{16} and boiling in the range of 130 °C to 290 °C (266 °F to 554 °F).]	294-799-5	91770-15-9	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-428-00-5	Kerosine (petroleum), solvent-refined sweetened; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from a petroleum stock by solvent refining and sweetening and boiling in the range of approximately 150 °C to 260 °C (302 °F to 500 °F).]	295-416-4	92045-36-8	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-429-00-0	Hydrocarbons, C_{9-16}, hydrotreated, dearomatized; Kerosine — unspecified; [A complex combination of hydrocarbons obtained as solvents which have been subjected to hydrotreatment in order to convert aromatics to naphthenes by catalytic hydrogenation.]	297-854-1	93763-35-0	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-430-00-6	Kerosine (petroleum), solvent-refined hydrodesulfurized; Kerosine — unspecified	307-033-2	97488-94-3	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-431-00-1	Distillates (petroleum), hydrodesulfurized full-range middle coker; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized coker distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_8 through C_{16} and boiling in the range of approximately 120 °C to 283 °C (248 °F to 541 °F).]	309-864-6	101316-58-9	Xn; R65	Xn R: 65 S: (2-)23-24-62		H

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-432-00-7	Solvent naphtha (petroleum), hydrodesulfurized heavy arom.; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by the catalytic hydrodesulfurization of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{13} and boiling in the range of approximately 180 °C to 240 °C (356 °F to 464 °F).]	309-882-4	101316-81-8	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-433-00-2	Solvent naphtha (petroleum), hydrodesulfurized medium; Kerosine — unspecified; [A complex combination of hydrocarbons obtained by the catalytic hydrodesulfurization of a petroleum fraction. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{13} and boiling in the range of approximately 175 °C to 220 °C (347 °F to 428 °F).]	309-884-5	101316-82-9	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-434-00-8	Kerosine (petroleum), hydrotreated; Kerosine — unspecified; [A complex combination of hydrocarbons obtained from the distillation of petroleum and subsequent hydrotreatment. It consists predominantly of alkanes, cycloalkanes and alkylbenzenes having carbon numbers predominantly in the range of C_{12} through C_{16} and boiling in the range of approximately 230 °C to 270 °C (446 °F to 518 °F).]	309-944-0	101631-19-0	Xn; R65	Xn R: 65 S: (2-)23-24-62		H
649-435-00-3	Distillates (petroleum), light catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F). It contains a relatively large proportion of bicyclic aromatic hydrocarbons.]	265-060-4	64741-59-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-436-00-9	Distillates (petroleum), intermediate catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{11} through C_{30} and boiling in the range of approximately 205 °C to 450 °C (401 °F to 842 °F). It contains a relatively large proportion of tricyclic aromatic hydrocarbons.]	265-062-5	64741-60-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-437-00-4	Distillates (petroleum), light hydrocracked; Cracked gasoil; [A complex combination of hydrocarbons from distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{18} and boiling in the range of approximately 160 °C to 320 °C (320 °F to 608 °F).]	265-078-2	64741-77-1	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
649-438-00-X	Distillates (petroleum), light thermal cracked; Cracked gasoil; [A complex combination of hydrocarbons from the distillation of the products from a thermal cracking process. It consists predominantly of unsaturated hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{22} and boiling in the range of approximately 160 °C to 370 °C (320 °F to 698 °F).]	265-084-5	64741-82-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-439-00-5	Distillates (petroleum), hydrodesulfurized light catalytic cracked; Cracked gasoil; [A complex combination of hydrocarbons obtained by treating light catalytic cracked distillates with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists of hydrocarbons having carbon numbers predominantly in the range of C_9 through C_{25} and boiling in the range of approximately 150 °C to 400 °C (302 °F to 752 °F). It contains a relatively large proportion of bicyclic aromatic hydrocarbons.]	269-781-5	68333-25-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-440-00-0	Distillates (petroleum), light steam-cracked naphtha; Cracked gasoil; [A complex combination of hydrocarbons from the multiple distillation of products from a steam cracking process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{10} through C_{18}.]	270-662-5	68475-80-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-441-00-6	Distillates (petroleum), cracked steam-cracked petroleum distillates; Cracked gasoil; [A complex combination of hydrocarbons produced by distilling cracked steam cracked distillate and/or its fractionation products. It consists of hydrocarbons having carbon numbers predominently in the range of C_{10} to low molecular weight polymers.]	270-727-8	68477-38-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-442-00-1	Gas oils (petroleum), steam-cracked; Cracked gasoil; [A complex combination of hydrocarbons produced by distillation of the products from a steam cracking process. It consists of hydrocarbons having carbon numbers predominantly greater than C_9 and boiling in the range of from approximately 205 °C to 400 °C (400 °F to 752 °F).]	271-260-2	68527-18-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-443-00-7	Distillates (petroleum), hydrodesulfurized thermal cracked middle; Cracked gasoil; [A complex combination of hydrocarbons obtained by fractionation from hydrodesulfurized themal cracker distillate stocks. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{11} to C_{25} and boiling in the range of approximately 205 °C to 400 °C (401 °F to 752 °F).]	285-505-6	85116-53-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-444-00-2	Gas oils (petroleum), thermal-cracked, hydrodesulfurized; Cracked gasoil	295-411-7	92045-29-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-445-00-8	Residues (petroleum), hydrogenated steam-cracked naphtha; Cracked gasoil; [A complex combination of hydrocarbons obtained as a residual fraction from the distillation of hydrotreated steam-cracked naphtha. It consists predominantly of hydrocarbons boiling in the range of approximately 200 °C to 350 °C (32 °F to 662 °F).]	295-514-7	92062-00-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-446-00-3	Residues (petroleum), steam-cracked naphtha distn.; Cracked gasoil; [A complex combination of hydrocarbons obtained as a column bottom from the separation of effluents from steam cracking naphtha at a high temperature. It boils in the range of approximately 147 °C to 300 °C (297 °F to 572 °F) and produces a finished oil having a viscosity of 18cSt at 50 °C.]	295-517-3	92062-04-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-447-00-9	Distillates (petroleum), light catalytic cracked, thermally degraded; Cracked gasoil; [A complex combination of hydrocarbons produced by the distillation of products from a catalytic cracking process which has been used as a heat transfer fluid. It consists predominantly of hydrocarbons boiling in the range of approximately 190 °C to 340 °C (374 °F to 644 °F). This stream is likely to contain organic sulfur compounds.]	295-991-1	92201-60-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-448-00-4	Residues (petroleum), steam-cracked heat-soaked naphtha; Cracked gasoil; [A complex combination of hydrocarbons obtained as residue from the distillation of steam cracked heat soaked naphtha and boiling in the range of approximately 150 °C to 350 °C (302 °F to 662 °F).]	297-905-8	93763-85-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-449-00-X	Hydrocarbons, C_{16-20}, solvent-dewaxed hydrocracked paraffinic distn. residue; Cracked gasoil; [A complex combination of hydrocarbons obtained by solvent dewaxing of a distillation residue from a hydrocracked paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{20} and boiling in the range of approximately 360 °C to 500 °C (680 °F to 932 °F). It produces a finished oil having a viscosity of 4,5 cSt at approximately 100 °C (212 °F).]	307-662-2	97675-88-2	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		H
649-450-00-5	Gas oils (petroleum), light vacuum, thermal-cracked hydrodesulfurized; Cracked gasoil; [A complex combination of hydrocarbons obtained by catalytic dehydrosulfurization of thermal-cracked light vacuum petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{20} and boiling in the range of approximately 270 °C to 370 °C (518 °F to 698 °F).]	308-278-8	97926-59-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-451-00-0	Distillates (petroleum), hydrodesulfurized middle coker; Cracked gasoil; [A complex combination of hydrocarbons by fractionation from hydrodesulfurised coker distillate stocks. Is consists of hydro-carbons having carbon numbers predominantly in the range of C_{12} through C_{21} and boiling in the range of approximately 200 °C to 360 °C (392 °F to 680 °F).]	309-865-1	101316-59-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H
649-452-00-6	Distillates (petroleum), heavy steam-cracked; Cracked gasoil; [A complex combination of hydrocarbons obtained by distillation of steam cracking heavy residues. It consists predominantly of highly alkylated heavy aromatic hydrocarbons boiling in the range of approximately 250 °C to 400 °C (482 °F to 752 °F).]	309-939-3	101631-14-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H

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649-453-00-1	Distillates (petroleum), heavy hydrocracked; Baseoil — unspecified; [A complex combination of hydrocarbons from the distillation of the products from a hydrocracking process. It consists predominantly of saturated hydrocarbons having carbon numbers in the range of C_{15}-C_{39} and boiling in the range of approximately 260 °C to 600 °C (500 °F to 1112 °F).]	265-077-7	64741-76-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-454-00-7	Distillates (petroleum), solvent-refined heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C).]	265-090-8	64741-88-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-455-00-2	Distillates (petroleum), solvent-refined light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-091-3	64741-89-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-456-00-8	Residual oils (petroleum), solvent deasphalted; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the solvent soluble fraction from C_3-C_4 solvent deasphalting of a residuum. It consists of hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-096-0	64741-95-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-457-00-3	Distillates (petroleum), solvent-refined heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt a 40 °C). It contains relatively few normal paraffins.]	265-097-6	64741-96-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-458-00-9	Distillates (petroleum), solvent-refined light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as the raffinate from a solvent extraction process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-098-1	64741-97-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-459-00-4	Residual oils (petroleum,) solvent-refined; Baseoil — unspecified; [A complex combination by hydrocarbons obtained as the solvent insoluble fraction from solvent refining of a residuum using a polar organic solvent such as phenol or furfural. It consists of hydrocarbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-101-6	64742-01-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-460-00-X	Distillates (petroleum), clay-treated paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-137-2	64742-36-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-461-00-5	Distillates (petroleum), clay-treated light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-138-8	64742-37-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-462-00-0	Residual oils (petroleum), clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treatment of a residual oil with a natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydro-carbons having carbon numbers predominantly higher than C_{25} and boiling above approximately 400 °C (752 °F).]	265-143-5	64742-41-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-463-00-6	Distillates (petroleum), clay-treated heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-146-1	64742-44-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-464-00-1	Distillates (petroleum), clay-treated light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contacting or percolation process to remove the trace amounts of polar compounds and impurities present. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-147-7	64742-45-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-465-00-7	Distillates (petroleum), hydrotreated heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-155-0	64742-52-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-466-00-2	Distillates (petroleum), hydrotreated light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-156-6	64742-53-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-467-00-8	Distillates (petroleum), hydrotreated heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-157-1	64742-54-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-468-00-3	Distillates (petroleum), hydrotreated light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains a relatively large proportion of saturated hydrocarbons.]	265-158-7	64742-55-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-469-00-9	Distillates (petroleum), solvent-dewaxed light paraffinic; Baseoil — unspecified; [A complex comination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-159-2	64742-56-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-470-00-4	Residual oils (petroleum), hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst. It consists of hydrocarbons having carbon numbers predominantly greater than C_{25} and boiling above approximately 400 °C (752 °F).]	265-160-8	64742-57-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-471-00-X	Residual oils (petroleum), solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of long, branched chain hydrocarbons from a residual oil by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly greater than C_{25} and boiling above approximately 400 °C (752 °F).]	265-166-0	64742-62-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-472-00-5	Distillates (petroleum), solvent-dewaxed heavy naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of not less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-167-6	64742-63-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-473-00-0	Distillates (petroleum), solvent-dewaxed light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists of hydrocarbons having carbon numbers predominantly in the range C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-168-1	64742-64-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-474-00-6	Distillates (petroleum), solvent-dewaxed heavy paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removal of normal paraffins from a petroleum fraction by solvent crystallization. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity not less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-169-7	64742-65-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-475-00-1	Naphthenic oils (petroleum), catalytic dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-172-3	64742-68-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-476-00-7	Naphthenic oils (petroleum), catalytic dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-173-9	64742-69-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-477-00-2	Paraffin oils (petroleum), catalytic dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C).]	265-174-4	64742-70-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-478-00-8	Paraffin oils (petroleum), catalytic dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewxing process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C).]	265-176-5	64742-71-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-479-00-3	Naphthenic oils (petroleum), complex dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by removing straight chain paraffin hydrocarbons as a solid by treatment with an agent such as urea. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of at least 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-179-1	64742-75-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-480-00-9	Naphthenic oils (petroleum), complex dewaxed light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from a catalytic dewaxing process. It consists of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	265-180-7	64742-76-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-481-00-4	Lubricating oils (petroleum), C_{20-50}, hydrotreated neutral oil-based, high-viscosity; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil, heavy vacuum gas oil, and solvent deasphalted residual oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil having a viscosity of approximately 112cSt at 40 °C. It contains a relatively large proportion of saturated hydrocarbons.]	276-736-3	72623-85-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-482-00-X	Lubricating oils (petroleum), C_{15-30}, hydrotreated neutral oil-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil and heavy vacuum gas oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil having a viscosity of approximately 15cSt at 40 °C. It contains a relatively large proportion of saturated hydrocabons.]	276-737-9	72623-86-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-483-00-5	Lubricating oils (petroleum), C_{20-50}, hydrotreated neutral oil-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating light vacuum gas oil, heavy vacuum gas oil and solvent deasphalted residual oil with hydrogen in the presence of a catalyst in a two stage process with dewaxing being carried out between the two stages. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of approximately 32cSt at 40 °C. It contains a relatively large proportion of saturated hydrocarbons.]	276-738-4	72623-87-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-484-00-0	Lubricating oils; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from solvent extraction and dewaxing processes. It consists predominantly of saturated hydrocarbons having carbon numbers in the range C_{15} through C_{50}.]	278-012-2	74869-22-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-485-00-6	Distillates (petroleum), complex dewaxed heavy paraffinci; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by dewaxing heavy paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil with a viscosity of equal to or greater than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	292-613-7	90640-91-8	Carc. Cat. 2; R45	T R: 4 S: 53-45		H L
649-486-00-1	Distillates (petroleum), complex dewaxed light paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by dewaxing light paraffinic distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{12} through C_{30} and produces a finished oil with a viscosity of less than 100 SUS at 100 °F (19cSt at 40 °C). It contains relatively few normal paraffins.]	292-614-2	90640-92-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-487-00-7	Distillates (petroleum), solvent dewaxed heavy paraffinic, clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating dewaxed heavy paraffinic distillate with neutral or modified clay in either a contacting or percolation process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	292-616-3	90640-94-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-488-00-2	Hydrocarbons, C_{20-50}, solvent dewaxed heavy paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons produced by treating dewaxed heavy paraffinic distillate with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	292-617-9	90640-95-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-489-00-8	Distillates (petroleum), solvent dewaxed light paraffinic, clay-treated; Baseoil — unspecified; [A complex combination of hydrocarbons resulting from treatment of dewaxed light paraffinic distillate with natural or modified clay in either a contacting or percolation process. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}.]	292-618-4	90640-96-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-490-00-3	Distillates (petroleum), solvent dewaxed light paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons produced by treating a dewaxed light paraffinic distillate with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}.]	292-620-5	90640-97-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-491-00-9	Residual oils (petroleum), hydrotreated solvent dewaxed; Baseoil — unspecified	292-656-1	90669-74-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-492-00-4	Residual oils (petroleum), catalytic dewaxed; Baseoil — unspecified	294-843-3	91770-57-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-493-00-X	Distillates (petroleum), dewaxed heavy paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from an intensive treatment of dewaxed distillate by hydrogenation in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{25} through C_{39} and produces a finished oil with a viscosity of approximately 44 cSt at 50 °C.]	295-300-3	91995-39-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-494-00-5	Distillates (petroleum), dewaxed light paraffinic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from an intensive treatment of dewaxed distillate by hydrogenation in the presence of a catalyst. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C_{21} through C_{29} and produces a finished oil with a viscosity of approximately 13 cSt at 50 °C.]	295-301-9	91995-40-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-495-00-0	Distillates (petroleum), hydrocracked solvent-refined, dewaxed; Baseoil — unspecified; [A complex combination of liquid hydrocarbons obtained by recrystallization of dewaxed hydrocracked solvent-refined petroleum distillates.]	295-306-6	91995-45-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-496-00-6	Distillates (petroleum), solvent-refined light naphthenic, hydrotreated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treating a petroleum fraction with hydrogen in the presence of a catalyst and removing the aromatic hydrocarbons by solvent extraction. It consists predominantly of naphthenic hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30} and produces a finished oil with a viscosity of between 13-15cSt at 40 °C.]	295-316-0	91995-54-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-497-00-1	Lubricating oils (petroleum), C_{17-35}, solvent-extd., dewaxed, hydrotreated; Baseoil — unspecified	295-423-2	92045-42-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-498-00-7	Lubricating oils (petroleum), hydrocracked nonarom. solvent-deparaffined; Baseoil — unspecified	295-424-8	92045-43-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-499-00-2	Residual oils (petroleum), hydrocracked acid-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons produced by solvent removal of paraffins from the residue of the distillation of acid-treated, hydrocracked heavy paraffins and boiling approximately above 380 °C (716 °F).]	295-499-7	92061-86-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-500-00-6	Paraffin oils (petroleum), solvent-refined dewaxed heavy; Baseoil — unspecified; [A complex combination of hydrocarbons obtained from sulfur-containing paraffinic crude oil. It consists predominantly of a solvent refined deparaffinated lubricating oil with a viscosity of 65cSt at 50 °C.]	295-810-6	92129-09-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-501-00-1	Lubricating oils (petroleum), base oils, paraffinic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by refining of crude oil. It consists predominantly of aromatics, naphthenics and paraffinics and produces a finished oil with a viscosity of 120 SUS at 100 °F (23cSt at 40 °C).]	297-474-6	93572-43-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-502-00-7	Hydrocarbons, hydrocracked paraffinic distn. residues, solvent-dewaxed; Baseoil — unspecified	297-857-8	93763-38-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-503-00-2	Hydrocarbons, C_{20-50}, residual oil hydrogenation vacuum distillate; Baseoil — unspecified	300-257-1	93924-61-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-504-00-8	Distillates (petroleum), solvent-refined hydrotreated heavy, hydrogenated; Baseoil — unspecified	305-588-5	94733-08-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-505-00-3	Distillates (petroleum), solvent-refined hydrocracked light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent dearomatization of the residue of hydrocracked petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{27} and boiling in the range of approximately 370 °C to 450 °C (698 °F to 842 °F).]	305-589-0	94733-09-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-506-00-9	Lubricating oils (petroleum), C_{18-40}, solvent-dewaxed hydrocracked distillate-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent deparaffination of the distillation residue from hydrocracked petroleum. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{40} and boiling in the range of approximately 370 °C to 550 °C (698 °F to 1022 °F).]	305-594-8	94733-15-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-507-00-4	Lubricating oils (petroleum), C_{18-40}, solvent-dewaxed hydrogenated raffinate-based; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent deparaffination of the hydrogenated raffinate obtained by solvent extraction of a hydrotreated petroleum distillate. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{40} and boiling in the range of approximately 370 °C to 550 °C (698 °F to 1022 °F).]	305-595-3	94733-16-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-508-00-X	Hydrocarbons, C_{13-30}, arom.-rich, solvent-extd. naphthenic distillate; Baseoil — unspecified	305-971-7	95371-04-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-509-00-5	Hydrocarbons, C_{16-32}, arom. rich, solvent-extd. naphthenic distillate; Baseoil — unspecified	305-972-2	95371-05-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-510-00-0	Hydrocarbons, C_{37-68}, dewaxed deasphalted hydrotreated vacuum distn. residues; Baseoil — unspecified	305-974-3	95371-07-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-511-00-6	Hydrocarbons, C_{37-65}, hydrotreated deasphalted vacuum distn. residues; Baseoil — unspecified	305-975-9	95371-08-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-512-00-1	Distillates (petroleum), hydrocracked solvent-refined light; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by the solvent treatment of a distillate from hydrocracked petroleum distillates. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{18} through C_{27} and boiling in the range of approximately 370 °C to 450 °C (698 °F to 842 °F.]	307-010-7	97488-73-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-513-00-7	Distillates (petroleum), solvent-refined hydrogenated heavy; Baseoil — unspecified; [A complex combination of hydrocarbons, obtained by the treatment of a hydrogenated petroleum distillate with a solvent. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{19} through C_{40} and boiling in the range of approximately 390 °C to 550 °C (734 °F to 1022 °F).]	307-011-2	97488-74-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-514-00-2	Lubricating oils (petroleum), C_{18-27}, hydrocracked solvent-dewaxed; Baseoil — unspecified	307-034-8	97488-95-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-515-00-8	Hydrocarbons, C_{17-30}, hydrotreated solvent-deasphalted atm. distn. residue, distn. lights; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the treatment of a solvent deasphalted short residue with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{30} and boiling in the range of approximately 300 °C to 400 °C (572 °F to 752 °F). It produces a finished oil having a viscosity of 4cSt at approximately 100 °C (212 °F).]	307-661-7	97675-87-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-516-00-3	Hydrocarbons, C_{17-40}, hydrotreated solvent-deasphalted distn. residue, vacuum distn. lights; Baseoil — unspecified; [A complex combination of hydrocarbons obtained as first runnings from the vacuum distillation of effluents from the catalytic hydrotreatment of a solvent deasphalted short residue having a viscosity of 8cSt at approximately 100 °C (212 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{40} and boiling in the range of approximately 300 °C to 500 °C (592 °F to 932 °F).]	307-755-8	97722-06-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-517-00-9	Hydrocarbons, C_{13-27}, solvent-extd. light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a viscosity of 9.5cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{27} and boiling in the range of approximately 240 °C to 400 °C (464 °F to 752 °F.]	307-758-4	97722-09-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-518-00-4	Hydrocarbons, C_{14-29}, solvent-extd. light naphthenic; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by extraction of the aromatics from a light naphthenic distillate having a viscosity of 16cSt at 40 °C (104 °F). It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{14} through C_{29} and boiling in the range of approximately 250 °C to 425 °C (482 °F to 797 °F).]	307-760-5	97722-10-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-519-00-X	Hydrocarbons, C_{27-42}, dearomatized; Baseoil — unspecified	308-131-8	97862-81-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-520-00-5	Hydrocarbons, C_{17-30}, hydrotreated distillates, distn. lights; Baseoil — unspecified	308-132-3	97862-82-3	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-521-00-0	Hydrocarbons, C_{27-45}, naphthenic vacuum distn.; Baseoil — unspecified	308-133-9	97862-83-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-522-00-6	Hydrocarbons, C_{27-45}, dearomatized; Baseoil — unspecified	308-287-7	97926-68-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-523-00-1	Hydrocarbons, C_{20-58}, hydrotreated; Baseoil — unspecified	308-289-8	97926-70-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-524-00-7	Hydrocarbons, C_{27-42}, naphthenic; Baseoil — unspecified	308-290-3	97926-71-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-525-00-2	Residual oils (petroleum), carbon-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by the treatment of solvent-dewaxed petroleum residual oils with activated charcoal for the removal of trace polar constituents and impurities.]	309-710-8	100684-37-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-526-00-8	Residual oils (petroleum), clay-treated solvent-dewaxed; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by treatment of solvent-dewaxed petroleum residual oils with bleaching earth for the removal of trace polar constituents and impurities.]	309-711-3	100684-38-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-527-00-3	Lubricating oils (petroleum), $C_{>25}$, solvent-extd., deasphalted, dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of vacuum distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly greater than C_{25} and produces a finished oil with a viscosity in the order of 32cSt to 37cSt at 100 °C (212 °F).]	309-874-0	101316-69-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-528-00-9	Lubricating oils (petroleum), C_{17-32}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{32} and produced a finished oil with a viscosity in the order of 17cSt to 23cSt at 40 °C (104 °F.]	309-875-6	101316-70-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-529-00-4	Lubricating oils (petroleum), C_{20-35}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{35} and produces a finished oil with a viscosity in the order of 37cSt to 44cSt at 40 °C (104 °F).]	309-876-1	101316-71-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-530-00-X	Lubricating oils (petroleum), C_{24-50}, solvent-extd., dewaxed, hydrogenated; Baseoil — unspecified; [A complex combination of hydrocarbons obtained by solvent extraction and hydrogenation of atmospheric distillation residues. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{24} through C_{50} and produces a finished oil with a viscosity in the order of 16cSt to 75cSt at 40 °C (104 °F).]	309-877-7	101316-72-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-531-00-5	Extracts (petroleum), heavy naphthenic distillate solvent, arom. conc.; Distillate aromatic extract (treated); [An aromatic concentrate produced by adding water to heavy naphthenic distillate solvent extract and extraction solvent.]	272-175-3	68783-00-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-532-00-0	Extracts (petroleum), solvent-refined heavy paraffinic distillate solvent; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from the re-extraction of solvent-refined heavy paraffinic distillate. It consists of saturated and aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	272-180-0	68783-04-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-533-00-6	Extracts (petroleum), heavy paraffinic distillates, solvent-deasphalted; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from a solvent extraction of heavy paraffinic distillate.]	272-342-0	68814-89-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-534-00-1	Extracts (petroleum), heavy naphthenic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by treating a heavy naphthenic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50} and produces a finished oil of at least 19cSt at 40 °C (100 SUS at 100 °F).]	292-631-5	90641-07-9	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-535-00-7	Extracts (petroleum), heavy paraffinic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons produced by treating a heavy paraffinic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{21} through C_{33} and boiling in the range of approximately 350 °C to 480 °C (662 °F to 896 °F).	292-632-0	90641-08-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

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649-536-00-2	Extracts (petroleum), light paraffinic distillate solvent, hydrotreated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons produced by treating a light paraffinic distillate solvent extract with hydrogen in the presence of a catalyst. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{17} through C_{26} and boiling in the range of approximately 280 °C to 400 °C (536 °F to 752 °F).]	292-633-6	90641-09-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-537-00-8	Extracts (petroleum), hydrotreated light paraffinic distillate solvent; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as the extract from solvent extraction of intermediate paraffinic top solvent distillate that is treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{36}.]	295-335-4	91995-73-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-538-00-3	Extracts (petroleum), light naphthenic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by treating the extract, obtained from a solvent extraction process, with hydrogen in the presence of a catalyst under conditions primarily to remove sulfur compounds. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{30}. This stream is likely to contain 5 wt.% or more of 4- to 6-membered condensed ring aromatic hydrocarbons.]	295-338-0	91995-75-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-539-00-9	Extracts (petroleum), light paraffinic distillate solvent, acid-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction of the distillation of an extract from the solvent extraction of light paraffinic top petroleum distillates that is subjected to a sulfuric acid refining. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	295-339-6	91995-76-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-540-00-4	Extracts (petroleum), light paraffinic distillate solvent, hydro-desulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of a light paraffin distillate and treated with hydrogen to convert the organic sulfur to hydrogen sulfide which is eliminated. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{40} and produces a finished oil with a viscosity of greater than 10cSt at 40 °C.]	295-340-1	91995-77-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-541-00-X	Extracts (petroleum), light vacuum gas oil solvent, hydro-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons, obtained by solvent extraction from light vacuum petroleum gas oils and treated with hydrogen in the presence of a catalyst. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	295-342-2	91995-79-8	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-542-00-5	Extracts (petroleum), heavy paraffinic distillate solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons resulting from treatment of a petroleum fraction with natural or modified clay in either a contact or percolation process to remove the trace amounts of polar compounds and impurities present. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}. This stream is likely to contain 5 wt.% or more 4-6 membered ring aromatic hydrocarbons.]	296-437-1	92704-08-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-543-00-0	Extracts (petroleum), heavy naphthenic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained from a petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{50} and produces a finished oil with a viscosity of greater than 19cSt at 40 °C.]	297-827-4	93763-10-1	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-544-00-6	Extracts (petroleum), solvent-dewaxed heavy paraffinic distillate solvent, hydrodesulfurized; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained from a solvent dewaxed petroleum stock by treating with hydrogen to convert organic sulfur to hydrogen sulfide which is removed. It consists predominantly of hydrocarbons having carbon numbers predominantly in the range of C_{15} through C_{50} and produces a finished oil with a viscosity of greater than 19cSt at 40 °C.]	297-829-5	93763-11-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-545-00-1	Extracts (petroleum), light paraffinic distillate solvent, carbon-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction from distillation of an extract recovered by solvent extraction of light paraffinic top petroleum distillate treated with activated charcoal to remove traces of polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	309-672-2	100684-02-4	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-546-00-7	Extracts (petroleum), light paraffinic distillate solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained as a fraction from distillation of an extract recovered by solvent extraction of light paraffinic top petroleum distillates treated with bleaching earth to remove traces of polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{16} through C_{32}.]	309-673-8	100684-03-5	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-547-00-2	Extracts (petroleum), light vacuum, gas oil solvent, carbon-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of light vacuum petroleum gas oil treated with activated charcoal for the removal of trace polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	309-674-3	100684-04-6	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-548-00-8	Extracts (petroleum), light vacuum gas oil solvent, clay-treated; Distillate aromatic extract (treated); [A complex combination of hydrocarbons obtained by solvent extraction of light vacuum petroleum gas oils treated with bleaching earth for removal of trace polar constituents and impurities. It consists predominantly of aromatic hydrocarbons having carbon numbers predominantly in the range of C_{13} through C_{30}.]	309-675-9	100684-05-7	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
649-549-00-3	Foots oil (petroleum); Foots oil; [A complex combination of hydrocarbons obtained as the oil fraction from a solvent deoiling or a wax sweating process. It consists predominantly of branched chain hydrocarbons having carbon numbers predominantly in the range of C_{20} through C_{50}.]	265-171-8	64742-67-2	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
649-550-00-9	Foots oil (petroleum), hydrotreated; Foots oil	295-394-6	92045-12-0	Carc. Cat. 2; R45	T R: 45 S: 53-45		H L
650-002-00-6	turpentine, oil	232-350-7	8006-64-2	R10 Xn; R20/21/22-65 Xi; R36/38 R43 N; R51-53	Xn; N R: 10-20/21/22-36/38-43-51/53-65 S: (2-)36/37-46-61-62		
650-003-00-1	fenson (ISO); 4-chlorophenyl benzenesulphonate;	201-274-6	80-38-6	Xn; R22 Xi; R36 N; R51-53	Xn; N R: 22-36-51/53 S: (2-)24-26-61		
650-004-00-7	norbormide (ISO); 5-(α-hydroxy-α-2-pyridylbenzyl)-7-(α-2-pyridylbenzylidene) bicyclo [2.2.1] hept-5-ene-2,3-dicarboximide	213-589-6	991-42-4	Xn; R22	Xn R: 22 S: (2-)		
650-005-00-2	(2*R*,6a*S*,12a*S*)-1,2,6,6a,12,12a-hexahydro-2-isopropenyl-8,9-dimethoxychromeno[3,4-b]furo[2,3-h]chromen-6-one, rotenone	201-501-9	83-79-4	T; R25 Xi; R36/37/38 N; R50-53	T; N R: 25-36/37/38-50/53 S: (1/2-)22-24/25-36-45-60-61		
650-006-00-8	benquinox (ISO); *p*-benzoquinone 1-benzoylhydrazone 4-oxime	207-807-9	495-73-8	T; R25 Xn; R21	T R: 21-25 S: (1/2-)36/37-45		
650-007-00-3	chlordimeform (ISO); N_2-(4-chloro-*o*-tolyl)-N_1,N_1-dimethylformamidine	228-200-5	6164-98-3	Carc. Cat. 3; R40 Xn; R21/22 N; R50-53	Xn; N R: 21/22-40-50/53 S: (2-)22-36/37-60-61		
650-008-00-9	drazoxolon (ISO); 4-(2-chlorophenylhydrazone)-3-methyl-5-isoxazolone	227-197-8	5707-69-7	T; R25 N; R50-53	T; N R: 25-50/53 S: (1/2-)22-24-36/37-45-60-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
650-009-00-4	chlordimeform hydrochloride; *N'*-(4-chloro-*o*-tolyl)-*N*,*N*-dimethylformamidine monohydrochloride; N^2-(4-chloro-*o*-tolyl)-N^1,N^1-dimethylformamidine hydorchloride	243-269-1	19750-95-9	Carc. Cat. 3; R40 Xn; R22 N; R50-53	Xn; N R: 22-40-50/53 S: (2-)22-36/37-60-61		
650-010-00-X	benzyl violet 4B; α-[4-(4-dimethylamino-α-{4-[ethyl(3-sodiosulphonatobenzyl)amino] phenyl}benzylidene)cyclohexa-2,5-dienylidene(ethyl)ammonio]toluene-3-sulphonate	216-901-9	1694-09-3	Carc. Cat. 3; R40	Xn R: 40 S: (2-)36/37		
650-012-00-0	erionite	—	12510-42-8	Carc. Cat. 1; R45	T R: 45 S: 53-45		
650-013-00-6	asbestos	— — — — — — —	12001-28-4 132207-32-0 12172-73-5 77536-66-4 77536-68-6 77536-67-5 12001-29-5	Carc. Cat. 1; R45 T; R48/23	T R: 45-48/23 S: 53-45		E
650-014-00-1	diethyl 2,4-dihydroxycyclodisiloxane-2,4-diylbis(trimethylene) diphosphonate, tetrasodium salt, reaction products with disodium metasilicate	401-770-4	—	C; R34 Xn; R22	C R: 22-34 S: (1/2-)26-36/37/39-45		
650-015-00-7	rosin; colophony	232-475-7 232-484-6 277-299-1	8050-09-7 8052-10-6 73138-82-6	R43	Xi R: 43 S: (2-)24-37		
650-016-00-2	Mineral wool, with the exception of those specified elsewhere in this Annex; [Man-made vitreous (silicate) fibres with random orientation with alkaline oxide and alkali earth oxide (Na_2O+K_2O+CaO +MgO+BaO) content greater than 18 % by weight]	—	—	Carc. Cat. 3; R40 Xi; R38	Xn R: 38-40 S: (2-)36/37		AQR
650-017-00-8	Refractory Ceramic Fibres; Special Purpose Fibres, with the exception of those specified elsewhere in this Annex; [Man-made vitreous (silicate) fibres with random orientation with alkaline oxide and alkali earth oxide (Na_2O+K_2O+CaO+ MgO+BaO) content less or equal to 18 % by weight]	—	—	Carc. Cat. 2; R49 Xi; R38	T R: 49-38 S: 53-45		A R

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
650-018-00-3	reaction product of: acetophenone, formaldehyde, cyclohexylamine, methanol and acetic acid	406-230-1	—	R10 Carc. Cat. 3; R40 C; R34 Xn; R20 R43 N; R50-53	C; N R: 10-20-34-40-43-50/53 S: (1/2-)26-36/37/39-45-60-61		
650-031-00-4	bis(4-hydroxy-*N*-methylanilinium) sulphate	200-237-1	55-55-0	Xn; R22-48/22 R43 N; R50-53	Xn; N R: 22-43-48/22-50/53 S: (2-)36/37-46-60-61		
650-032-00-X	cyproconazole (ISO); (2*RS*,3*RS*;2*RS*,3*SR*)-2-(4-chlorophenyl)-3-cyclopropyl-1-(1*H*-1,2,4-triazol-1-yl)butan-2-ol	—	94361-06-5	Repr. Cat. 3; R63 Xn; R22 N; R50-53	Xn; N R: 22-50/53-63 S: (2-)36/37-60-61		
650-033-00-5	(*S*)-α-cyano-3-phenoxybenzyl-(*S*)-2-(4-chlorophenyl)-3-methylbutyrate; esfenvalerate	—	66230-04-4	T; R23/25 R43 N; R50-53	T; N R: 23/25-43-50/53 S: (1/2-)24-36/37/39-45-60-61		
650-041-00-9	triasulfuron (ISO); 1-[2-(2-chloroethoxy)phenylsulfonyl]-3-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)urea	—	82097-50-5	N; R50-53	N R: 50/53 S: 60-61		
650-042-00-4	reaction product of: polyethylene-polyamine-(C_{16}-C_{18})-alkylamides with monothio-(C_2)-alkyl phosphonates	417-450-2	—	Xi; R36/38 R43 R52-53	Xi R: 36/38-43-52/53 S: (2-)24-26-37-61		
650-043-00-X	reaction product of: 3,5-bis-*tert*-butylsalicylic acid and aluminiumsulfate	420-310-3	—	Xn; R22 N; R50-53	Xn; N R: 22-50/53 S: (2-)22-56-60-61		
650-044-00-5	mixed linear and branched C_{14-15} alcohols ethoxylated, reaction product with epichlorohydrin	420-480-9	158570-99-1	Xi; R38 R43 N; R50-53	Xi; N R: 38-43-50/53 S: (2-)24-37-60-61		
650-045-00-0	reaction product of: 1,2,3-propanetricarboxylic acid, 2-hydroxy, diethyl ester, 1-propanol and zirconium tetra-n-propanolate	417-110-3	—	F; R11 Xi; R38-41 N; R51-53	F; Xi; N R: 11-38-41-51/53 S: (2-)9-16-26-37/39-61		

Index No	International Chemical Identification	EC No	CAS No	Classification	Labelling	Concentration Limits	Notes
650-046-00-6	di(tetramethylammonium)(29*H*,31*H*-phthalocyanin-*N*29, *N*30,*N*31,*N*32)disulfonamide disulfonate, cuprate(2-)complex, derivates	416-180-2	12222-04-7	Xn; R22-48/22 N; R51-53	Xn; N R: 22-48/22-51/53 S: (2-)22-36-61		
650-047-00-1	dibenzylphenylsulfonium hexafluoroantimonate	417-760-8	134164-24-2	T; R48/25 Xn; R22 Xi; R41 R43 N; R51-53	T; N R: 22-41-43-48/25-51/53 S: (1/2-)22-26-36/37/39-45-61		
650-048-00-7	reaction product of: borax, hydrogen peroxide, acetic acid anhydride and acetic acid	420-070-1	—	O; R7 Xn; R20/21/22 C; R35 N; R50	O; C; N R: 7-20/21/22-35-50 S: (1/2-)3/7-14-26-36/37/39-45-61		
650-049-00-2	2-alkoyloxyethyl hydrogen maleate, where alkoyl represents (by weight) 70 to 85 % unsaturated octadecoyl, 0.5 to 10 % saturated octadecoyl, and 2 to 18 % saturated hexadecoyl	417-960-5	—	Xi; R38-41 R43 N; R50-53	Xi; N R: 38-41-43-50/53 S: (2-)24-26-37/39-60-61		
650-050-00-8	reaction mass of: 1-methyl-3-hydroxypropyl 3,5-[1,1-dimethylethyl]-4-hydroxydihydro-cinnamate and/or 3-hydroxybutyl 3,5-[1,1-dimethylethyl]-4-hydroxydihydrocinnamate; 1,3-butanediol bis[3-(3'-(1,1-dimethylethyl)4'-hydroxy-phenyl)propionate] isomers; 1,3-butanediol bis[3-(3',5'-(1,1-dimethylethyl)-4'-hydroxyphenyl)propionate] isomers	423-600-8	—	N; R51-53	N R: 51/53 S: 61		
650-055-00-5	silver sodium zirconium hydrogenphosphate	422-570-3	155925-27-2	N; R50-53	N R: 50/53 S: 60-61		

ANNEX VII

Translation table from classification under Directive 67/548/EEC to classification under this Regulation

This Annex includes a table to assist translation of a classification made for a substance or a mixture under Directive 67/548/EEC or Directive 1999/45/EC, respectively, into the corresponding classification under this Regulation. Whenever data for the substance or mixture are available, an evaluation and classification shall be done in accordance with Articles 9 to13 of this Regulation.

1. Translation table

The codes used are introduced in Table 1.1 and section 1.1.2.2 of Annex VI.

Table 1.1

Translation between classification in accordance with Directive 67/548/EEC and this Regulation

Classification under Directive 67/548/EEC	Physical state of the substance when relevant	Classification under this Regulation		Note
		Hazard Class-and-Category	Hazard statement	
E; R2		No direct translation possible.		
E; R3		No direct translation possible.		
O; R7		Org. Perox. CD	H242	
		Org. Perox. EF	H242	
O; R8	gas	Ox. Gas 1	H270	
O; R8	liquid, solid	No direct translation possible.		
O; R9	liquid	Ox. Liq. 1	H271	
O; R9	solid	Ox. Sol. 1	H271	
R10	liquid	No direct translation possible. Correct translation of R10, liquid is: — Flam. Liq. 1, H224 if flashpoint < 23 °C and initial boiling point ≤ 35 °C — Flam. Liq. 2, H225 if flashpoint < 23 °C and initial boiling point > 35 °C — Flam. Liq. 3, H226 if flashpoint ≥ 23 °C		
F; R11	liquid	No direct translation possible. Correct translation of F; R11, liquid is: — Flam. Liq. 1, H224 if initial boiling point ≤ 35 °C — Flam. Liq. 2, H225 if initial boiling point > 35 °C		
F; R11	solid	No direct translation possible.		
F+; R12	gas	No direct translation possible. Correct translation of F+; R12, gaseous results either in Flam. Gas 1, H220 or Flam. Gas 2, H221.		
F+; R12	liquid	Flam. Liq. 1	H224	
F+; R12	liquid	Self-react. CD	H242	
		Self-react. EF	H242	
		Self-react. G	none	
F; R15		No translation possible.		
F; R17	liquid	Pyr. Liq. 1	H250	
F; R17	solid	Pyr. Sol. 1	H250	

Classification under Directive 67/548/EEC	Physical state of the substance when relevant	Classification under this Regulation		Note
		Hazard Class-and-Category	Hazard statement	
Xn; R20	gas	Acute Tox. 4	H332	(1)
Xn; R20	vapours	Acute Tox. 4	H332	(1)
Xn; R20	dust/mist	Acute Tox. 4	H332	
Xn; R21		Acute Tox. 4	H312	(1)
Xn; R22		Acute Tox. 4	H302	(1)
T; R23	gas	Acute Tox. 3	H331	(1)
T; R23	vapour	Acute Tox. 2	H330	
T; R23	dust/mist	Acute Tox. 3	H331	(1)
T; R24		Acute Tox. 3	H311	(1)
T; R25		Acute Tox. 3	H301	(1)
T+; R26	gas	Acute Tox. 2	H330	(1)
T+; R26	vapour	Acute Tox. 1	H330	
T+; R26	dust/mist	Acute Tox. 2	H330	(1)
T+; R27		Acute Tox. 1	H310	
T+; R28		Acute Tox. 2	H300	(1)
R33		STOT RE 2	H373	(3)
C; R34		Skin Corr. 1B	H314	(2)
C; R35		Skin Corr. 1A	H314	
Xi; R36		Eye Irrit. 2	H319	
Xi; R37		STOT SE 3	H335	
Xi; R38		Skin Irrit. 2	H315	
T; R39/23		STOT SE 1	H370	(3)
T; R39/24		STOT SE 1	H370	(3)
T; R39/25		STOT SE 1	H370	(3)
T+; R39/26		STOT SE 1	H370	(3)
T+; R39/27		STOT SE 1	H370	(3)
T+; R39/28		STOT SE 1	H370	(3)
Xi; R41		Eye Dam. 1	H318	
R42		Resp. Sens. 1	H334	
R43		Skin Sens. 1	H317	
Xn; R48/20		STOT RE 2	H373	(3)
Xn; R48/21		STOT RE 2	H373	(3)
Xn; R48/22		STOT RE 2	H373	(3)
T; R48/23		STOT RE 1	H372	(3)
T; R48/24		STOT RE 1	H372	(3)
T; R48/25		STOT RE 1	H372	(3)

Classification under Directive 67/548/EEC	Physical state of the substance when relevant	Classification under this Regulation		Note
		Hazard Class-and-Category	Hazard statement	
R64		Lact.	H362	
Xn; R65		Asp. Tox. 1	H304	
R67		STOT SE 3	H336	
Xn; R68/20		STOT SE 2	H371	(3)
Xn; R68/21		STOT SE 2	H371	(3)
Xn; R68/22		STOT SE 2	H371	(3)
Carc. Cat. 1; R45		Carc. 1A	H350	
Carc. Cat. 2; R45		Carc. 1B	H350	
Carc. Cat. 1; R49		Carc. 1A	H350i	
Carc. Cat. 2; R49		Carc. 1B	H350i	
Carc. Cat. 3; R40		Carc. 2	H351	
Muta. Cat. 2; R46		Muta. 1B	H340	
Muta. Cat. 3; R68		Muta. 2	H341	
Repr. Cat. 1; R60		Repr. 1A	H360F	(4)
Repr. Cat. 2; R60		Repr. 1B	H360F	(4)
Repr. Cat. 1; R61		Repr. 1A	H360D	(4)
Repr. Cat. 2; R61		Repr. 1B	H360D	(4)
Repr. Cat. 3; R62		Repr. 2	H361f	(4)
Repr. Cat. 3; R63		Repr. 2	H361d	(4)
Repr. Cat. 1; R60-61		Repr. 1A	H360FD	
Repr. Cat. 1; R60 Repr. Cat. 2; R61		Repr. 1A	H360FD	
Repr. Cat. 2; R60 Repr. Cat. 1; R61		Repr. 1A	H360FD	
Repr. Cat. 2; R60-61		Repr. 1B	H360FD	
Repr. Cat. 3; R62-63		Repr. 2	H361fd	
Repr. Cat. 1; R60 Repr. Cat. 3; R63		Repr. 1A	H360Fd	
Repr. Cat. 2; R60 Repr. Cat. 3; R63		Repr. 1B	H360Fd	
Repr. Cat. 1; R61 Repr. Cat. 3; R62		Repr. 1A	H360Df	
Repr. Cat. 2; R61 Repr. Cat. 3; R62		Repr. 1B	H360Df	
N; R50		Aquatic. Acute 1	H400	
N; R50-53		Aquatic Acute 1 Aquatic Chronic 1	H400 H410	
N; R51-53		Aquatic Chronic 2	H411	
R52-53		Aquatic Chronic 3	H412	

Classification under Directive 67/548/EEC	Physical state of the substance when relevant	Classification under this Regulation		Note
		Hazard Class-and-Category	Hazard statement	
R53		Aquatic Chronic 4	H413	
N; R59		Ozone	EUH059	

Note 1

For these classes it is possible to use the recommended minimum classification as defined in section 1.2.1.1 in Annex VI. Data or other information may be available to indicate that re-classification in a more severe category is appropriate.

Note 2

It is recommended to classify in Category 1B even if it also could be possible that 1C could be applicable for certain cases. Going back to original data, may not result in a possibility to distinguish between Category 1B or 1C, since the exposure period has normally been up to 4 hours according to Regulation (EC) No 440/2008. However, for the future, when data are derived from tests following a sequential approach as foreseen in the Regulation (EC) No 440/2008, Category 1C should be considered.

Note 3

The route of exposure could be added to the hazard statement if it is conclusively proven that no other routes of exposure cause the hazard.

Note 4

Hazard statements H360 and H361 indicate a general concern for both the reproductive properties related to fertility and developmental effects; 'May damage/Suspected of damaging fertility or the unborn child'. According to the classification criteria (Annex I, section 3.7) the general hazard statement can be replaced by the hazard statement indicating only the property of concern, in case either fertility or developmental effects are proven to be not relevant.

Table 1.2

Translation between risk phrases assigned under Directive 67/548/EEC and supplementary labelling requirements under this Regulation

Directive 67/548/EEC	This Regulation
R1	EUH001
R6	EUH006
R14	EUH014
R18	EUH018
R19	EUH019
R44	EUH044
R29	EUH029
R31	EUH031
R32	EUH032
R66	EUH066
R39-41	EUH070