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MINISTRY OF HEALTH OF THE RUSSIAN FEDERATION

**CHIEF STATE SANITARY INSPECTOR
OF THE RUSSIAN FEDERATION**

**RESOLUTION
No. 24 dated September 26, 2001**

ON ENACTMENT OF SANITARY RULES

(as amended by Resolution No. 20 of the Chief State Sanitary Inspector of the Russian Federation dated April 7, 2009,
Amendment No. 2 approved by Resolution No. 10 of the Chief State Sanitary Inspector of the Russian Federation dated February 25, 2010, as amended by Amendment No.3 approved by Resolution No. 74 of the Chief State Sanitary Inspector of the Russian Federation dated June 28, 2010)

On the basis of Federal Law No. 52-FZ "On Sanitary and Epidemiological Well-Being of Population" dated March 30, 1999 <1> and the Regulation on State Sanitary and Epidemiological Control <2> approved by Resolution No. 554 of the Government of the Russian Federation dated July 24, 2000, I hereby order:

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- <1> Collection of laws and regulations of the Russian Federation No. 14, 1999, Article 1650.
 - <2> Collection of laws and regulations of the Russian Federation No. 31, 2000, Article 3295.

1. To put in force the sanitary and epidemiological rules and regulations "Drinking Water. Hygienic Requirements for Quality of Water in the Centralized Drinking Water Supply Systems. Quality Control. SanPiN 2.1.4.1074-01" approved by the Chief State Sanitary Inspector of the Russian Federation on September 26, 2001, from January 1, 2002.

G. G. ONISHCHENKO

Approved by
the Chief State
Sanitary Inspector
of the Russian Federation
on September 26, 2001

2.1.4. DRINKING WATER AND WATER SUPPLY FOR THE POPULATED AREAS

Amendment No. 3 to these Sanitary Rules, which contains a separate chapter "Hygienic Requirements for Safety of Materials, Chemical Reagents and Equipment Used for Water Purification and Treatment", was approved by Resolution No. 74 of the Chief State Sanitary Inspector of the Russian Federation dated June 28, 2010.

The Amendment to these Sanitary Rules which contains a separate chapter "Hygienic Requirements for Hot Water Supply Systems Safety Assurance" brought into force from September 1, 2009 was approved by Resolution No. 20 of the Chief State Sanitary Inspector of the Russian Federation dated April 7, 2009.

**DRINKING WATER. HYGIENIC REQUIREMENTS FOR QUALITY OF WATER
IN THE CENTRALIZED DRINKING WATER SUPPLY SYSTEMS.
QUALITY CONTROL. HYGIENIC REQUIREMENTS FOR HOT WATER
SUPPLY SYSTEMS SAFETY ASSURANCE**

SANITARY AND EPIDEMIOLOGICAL RULES AND REGULATIONS
SanPiN 2.1.4.1074-01

(as amended by Resolution No. 20 of the Chief State Sanitary Inspector of the Russian Federation dated April 7, 2009,
Amendment No.2 approved by Resolution No. 10 of the Chief State Sanitary Inspector of the Russian Federation dated February 25, 2010,
as amended by Amendment No.3 approved by Resolution No. 74 of the Chief State Sanitary Inspector of the Russian Federation dated June 28, 2010)

1. Scope of Application

1.1. Sanitary and epidemiological rules and regulations "Drinking Water. Hygienic Requirements for Quality of Water in the Centralized Drinking Water Supply Systems. Quality Control" (hereinafter referred to as the "Sanitary Rules") set forth hygienic requirements for the quality of the drinking water as well as regulations for control over the quality of water processed and delivered by the populated areas centralized drinking water supply systems (hereinafter referred to as the "Water supply systems").

Note:

The Regulations on the State Sanitary and Epidemiological Service of the Russian Federation approved by Resolution No. 554 of the Government of the Russian Federation dated July 24, 2000 ceased to be in force due to enactment of Resolution No. 569 of the Government of the Russian Federation dated September 15, 2005.

1.2. These Sanitary Rules have been developed on the basis of Federal Law "On Sanitary and Epidemiological Well-Being of Population", Fundamentals of Legislation of the Russian Federation on Public Health Care <1>, Regulations on State Sanitary and Epidemiological Control and Regulations on State Sanitary and Epidemiological Service of the Russian Federation <2>.

<1> Bulletin of the Congress of People's Deputies of the Russian Federation and the Supreme Council of the Russian Federation No. 33, 1993, Article 1318.

<2> Collection of laws and regulations of the Russian Federation No. 31, 2000, Article 3295.

1.3. The Sanitary Rules are intended for private entrepreneurs and legal entities which activity is related to design, construction and operation of the water supply systems and provision of the population with drinking water as well as for agencies and organizations which carry out sanitary and epidemiological supervision and control.

1.4. The Sanitary Rules apply to the water delivered by the water supply systems and intended for population consumption for drinking and household needs, for using in food raw materials processing and production of food, storage and trading thereof as well as for manufacturing of products requiring water of drinking quality.

1.5. Hygienic requirements for the quality of drinking water in case of non-centralized water supply, for the quality of drinking water processed by autonomous water supply systems, individual water treatment facilities as well as of water sold to the population in bottles and containers shall be set forth by other sanitary rules and regulations.

2. General Provisions

2.1. The requirements of these Sanitary Rules shall be complied with in development of state standards, construction rules and regulations governing drinking water supply to the population, design and technical documents for the water supply systems as well as during construction and operation of the water supply systems.

2.2. The quality of drinking water delivered by the water supply system shall conform to the requirements of these Sanitary Rules.

2.3. The parameters characterizing regional peculiarities of the drinking water chemical composition shall be set individually for each water supply system in accordance with the rules specified in Supplement 1.

2.4. Based on the requirements of these Sanitary Rules a private entrepreneur or a legal entity operating the water supply system shall develop a work program for the water quality production control (hereinafter referred to as the "Work Program") in accordance with the rules specified in Supplement 1. The Work Program shall be approved by the municipal or regional State Sanitary and Epidemiological Supervision Center (hereinafter referred to as the "State Sanitary and Epidemiological Supervision Center") and shall be adopted for the corresponding territory within the specified procedure.

2.5. Subject to emergency situations or technical violations at objects and facilities of the water supply system that result or can result in decrease of the drinking water quality and conditions of the water supply to the population, a private entrepreneur or a legal entity operating the water supply system shall immediately take measures to eliminate the said situations or violations and inform the State Sanitary and Epidemiological Supervision Center thereof.

A private entrepreneur or a legal entity exercising production control of the drinking water quality shall also immediately inform the State Sanitary and Epidemiological Supervision Center about each result of water sample laboratory analysis which fails to conform with the hygienic standards.

2.6. In case of natural phenomena which cannot be foreseen in advance or in case of emergency situations which cannot be eliminated immediately, there might be temporary divergences from the hygienic standards of the drinking water quality allowed only for the chemical composition parameters affecting the organoleptic attributes.

2.6.1. Divergences from the hygienic standards are allowed only subject to concurrence of the following conditions:

- drinking water cannot be supplied to the population by any other means;
- the maximum permissible divergences from the hygienic standards approved for a limited period of time by the State Sanitary and Epidemiological Supervision Center are complied with;
- the maximum time limits are complied with;
- there is no risk for health of the population during the period of divergences;
- the population is informed about the said divergences and the duration thereof, about absence of risks for health as well as about recommendations on the use of drinking water.

2.6.2. The decision on temporary divergence from the hygienic standards of the drinking water quality shall be made in accordance with the legislation of the Russian Federation.

2.6.3. An action plan for provision of the water quality conforming to the hygienic standards (including a calendar work plan, deadlines and funds available) shall be adopted upon taking the decision on temporary divergence from the hygienic standards.

2.7. Drinking water supply to the population shall be prohibited or its use shall be suspended in the following cases:

- the reasons for deterioration of the drinking water quality are not eliminated within the specified period of temporary divergences from the hygienic standards;
- the water supply system fails to produce and supply to the population drinking water the quality of which conforms with the requirements of these Sanitary Rules thus creating real hazard to the population health.

2.7.1. The decision on prohibition or suspension of the population's use of the drinking water from the specific water supply system shall be made by the local authorities following a resolution of the Chief State Sanitary Inspector of the corresponding territory based on the assessment of hazard and risks for the population's health arising due to both further consumption of water that does not conform to the hygienic standards and prohibition or suspension of the said water use for drinking and household needs.

2.7.2. In case of a decision on prohibition or suspension of the drinking water use, the organizations ensuring operation of the water supply system shall (subject to the agreement with the State Sanitary and Epidemiological Supervision Center) develop and carry out actions aimed at detection and elimination of the reasons for the water quality deterioration as well as at provision of the population with drinking water conforming with the requirements of the Sanitary Rules.

2.7.3. The population shall be informed about the decision on prohibition or suspension of the drinking water use, the water quality, the actions carried out as well as about recommendations on acting in this situation within the specified procedure.

3. Hygienic Requirements and Standards for Drinking Water Quality

3.1. Drinking water shall be safe epidemiologically and radiologically, have benign chemical composition and good organoleptic attributes.

3.2. The drinking water quality shall conform to the hygienic standards before supply to the distribution facilities as well as at the water supply points of the external and internal water supply network.

3.3. Epidemiological safety of drinking water is defined by its conformity to the standards for microbiological and parasitological parameters provided in Table 1.

Table 1

| Parameters | Measurement units | Standards |
|---|---|-----------|
| Thermotolerant Coliform bacteria | Amount of bacteria per 100 ml <1> | None |
| General Coliform bacteria <2> | Amount of bacteria per 100 ml <1> | None |
| Total bacterial count <2> | Amount of colony-forming units per 1 ml | max 50 |
| Coliphages <3> | Amount of the plaque-forming units (PFU) per 100 ml | None |
| Sulfite-reducing Clostridium spores <4> | Amount of spores per 20 ml | None |
| Lambliа cysts <3> | Amount of cysts per 50 liters | None |

Notes.

<1> Triple analysis of 100 ml of the water samples is conducted for the determination.

<2> The exceedance of the standard is not allowed for 95% of samples collected at the water supply points of the external and internal water supply network within 12 months in case of at least 100 analyzed samples per year.

<3> The determination is conducted only in the water supply systems where the water is taken from the surface sources before its supply to the distribution network.

<4> The determination is conducted within water technology effectiveness assessment.

3.3.1. Determination of thermotolerant coliform bacteria, general coliform bacteria, the total bacterial count and coliphages is conducted at each sample when analysing the microbiological parameters of the drinking water quality.

3.3.2. Should thermotolerant coliform bacteria and (or) general coliform bacteria and (or) coliphages be detected in a drinking water sample, these are to be determined using immediately repeatedly collected water samples. In such cases the determination of chlorides, ammonia nitrogen, nitrates and nitrites shall be simultaneously conducted in order to find the reasons for contamination.

3.3.3. Should general coliform bacteria exceed 2 per 100 ml, and (or) thermotolerant coliform bacteria and (or) coliphages be found in the repeatedly collected water samples, the water samples analysis to determine pathogenic bacteria of the coliform bacterium group and (or) enteroviruses shall be conducted.

3.3.4. Analyses of drinking water for pathogenic bacteria of the coliform bacterium group and enteroviruses are also conducted on epidemiological indications following the decision of the State Sanitary and Epidemiological Supervision Center.

3.3.5. Analyses of water for pathogenic microorganisms can be conducted only in laboratories having sanitary-epidemiological conclusion on the compliance of the work conditions with the sanitary rules and a license for activities related to the use of infectious agents.

3.4. Safety of drinking water as to its chemical composition shall be determined by its conformity to the standards on:

3.4.1. the composite index and hazardous chemical content of substances which are the most frequent in the natural waters in the territory of the Russian Federation as well as of globally distributed man-made substances (Table 2);

3.4.2. the content of harmful chemical substances discharged to and formed in water during its processing in the water supply system (Table 3);

3.4.3. the content of harmful chemical substances discharged to the water supply sources as a result of anthropogenic activities (Supplement 2).

Table 2

| Parameters | Units of measurement | Standards (maximum allowable concentration (MAC)), up to | Hazard index <1> | Class of hazard |
|--|----------------------|--|------------------|-----------------|
| Composite indices | | | | |
| Hydrogen ion concentration | units, pH | in the range of 6–9 | | |
| Total salt content (solid residue) | mg/l | 1000 (1500) <2> | | |
| Total water hardness | mg-equivalents/l | 7.0 (10) <2> | | |
| Permanganate demand | mg/l | 5.0 | | |
| Petrochemicals, cumulative | mg/l | 0.1 | | |
| Surface active substances (SAS), anionic | mg/l | 0.5 | | |
| Phenol index | mg/l | 0.25 | | |
| Inorganic Substances | | | | |
| Aluminum (AL ³⁺) | mg/l | 0.5 | s.-t. | 2 |
| Barium (Ba ²⁺) | mg/l | 0.1 | s.-t. | 2 |
| Beryllium (Be ²⁺) | - " | 0.0002 | - " | 1 |
| Boron (B, cumulative) | - " | 0.5 | - " | 2 |
| Ferrum (Fe, cumulative) | - " | 0.3 (1.0) <2> | org. | 3 |
| Cadmium (Cd, cumulative) | - " | 0.001 | s.-t. | 2 |

| | | | | |
|--|------|---------------|-------|---|
| Manganese (Mn, cumulative) | - " | 0.1 (0.5) <2> | org. | 3 |
| Cuprum (B, cumulative) | - " | 1.0 | - " | 3 |
| Molybdenum (Mo, cumulative) | - " | 0.25 | s.-t. | 2 |
| Arsenium (As, cumulative) | - " | 0.05 | s.-t. | 2 |
| Nickel (Ni, cumulative) | mg/l | 0.1 | s.-t. | 3 |
| Nitrates (as to NO ₃) | - " | 45 | s.-t. | 3 |
| Mercury (Hg, cumulative) | - " | 0.0005 | s.-t. | 1 |
| Plumbum (Pb, cumulative) | - " | 0.03 | - " | 2 |
| Selenium (Se, cumulative) | - " | 0.01 | - " | 2 |
| Strontium (Sr ²⁺) | - " | 7.0 | - " | 2 |
| Sulphates (SO ₄ ²⁻) | - " | 500 | org. | 4 |
| Fluorides (F ⁻) | | | | |
| for climatic areas | | | | |
| - I и II | - " | 1.5 | s.-t. | 2 |
| - III | - " | 1.2 | | 2 |
| Chlorides (Cl ⁻) | - " | 350 | org. | 4 |
| Chromium (Cr ⁶⁺) | - " | 0.05 | s.-t. | 3 |
| Cyanides (CN ⁻) | - " | 0.035 | - " | 2 |
| Zink (Zn ²⁺) | - " | 5.0 | org. | 3 |
| Organic Substances | | | | |

| | | | | | |
|----------------------|------|-----------|-------|---|--|
| gamma HCCH (lindane) | - "- | 0.002 <3> | s.-t. | 1 | |
| DDT (total isomers) | - "- | 0.002 <3> | - "- | 2 | |
| 2,4-D | - "- | 0.03 <3> | - "- | 2 | |

Notes.

<1> The limiting harmful index according to which the standard is set: "s.-t." — sanitary and toxicological, "org." — organoleptic.

<2> The figure in brackets can be set according to the Resolution of the Chief State Sanitary Inspector of the corresponding territory for a specific water supply system based on the sanitary and epidemiological situation in the residential place and the water treatment technology applied.

<3> The standards are adopted in accordance with the WHO recommendations.

Table 3

| Parameters | Units of measurement | Standards (maximum allowable concentration (MAC)), up to | Hazard index | Class of hazard | |
|---|----------------------|--|--------------|-----------------|--|
| Chlorum <1> | | | | | |
| - residual free | mg/l | within 0.3 - 0.5 | org. | 3 | |
| - residual fixed | - "- | within 0.8 - 1.2 | - "- | 3 | |
| Chloroform (at water chlorination) | - "- | 0.2 <2> | s.-t. | 2 | |
| Residual ozone <3> | - "- | 0.3 | org. | | |
| Formaldehyde (at water ozone treatment) | - "- | 0.05 | s.-t. | 2 | |
| Polyacrylamide | - "- | 2.0 | - "- | 2 | |
| Activated silicic acid (per Si) | - "- | 10 | - "- | 2 | |
| Polyphosphates (per PO ₄ ³⁻) | - "- | 3.5 | org. | 3 | |
| Residual amount of aluminum and ferrum bearing coagulants | - "- | See parameters "Aluminum", "Ferrum" in Table 2 | | | |

Notes.

<1> At sterilization of water with the free chlorine the time of its contact with water shall be at least 30 minutes, with the fixed chlorine—at least 60 minutes.

Control over the content of the residual chlorine is conducted before the water supply into the distribution network.

Should free and fixed chlorine simultaneously be in water, their cumulative concentration shall not exceed 1.2 mg/l.

In the exceptional cases subject to agreement with the State Sanitary and Epidemiological Supervision Center the increased concentration of chlorine in the drinking water can be allowed.

<3> Standard adopted in accordance with the WHO recommendations.

<3> Control over the content of the residual ozone is conducted after the mixing chamber should the contact time be at least 12 minutes.

3.4.4. Should several chemical substances of the 1st and 2nd hazard class standardized as per their sanitary and toxicological health hazard characteristic be found in the drinking water, the cumulative ratio of each substance concentration in the water to its MAC shall not exceed 1. The calculation is done according to the following formula:

$$\frac{C_1 \text{ fact}}{C_1 \text{ adm}} + \frac{C_2 \text{ fact}}{C_2 \text{ adm}} + \dots + \frac{C_n \text{ fact}}{C_n \text{ adm}} \leq 1,$$

where C_1, C_2, C_n indicate the concentration of individual chemical substances of the 1st and 2nd hazard class: fact. (factual) and adm. (admissible).

3.5. Propitious organoleptic attributes of water are determined by its conformance to the standards specified in Table 4 as well as to standards for the content of substances influencing the organoleptic attributes of water provided in Table 2 and Table 3 as well as in Supplement 2.

Table 4

| Parameters | Units of measurement | Standards max |
|-----------------------------------|---|------------------------------|
| Odor | points | 2 |
| Flavour | - "— | 2 |
| Colority | degrees | 20 (35) <1> |
| Suspended-materials concentration | FTU (Formazin Turbidity Units) or mg/l (per kaoline) | 2,6 (3,5) <1> 1,5 (2) <1> |

Note.

<1> The figure in brackets can be set following the Resolution of Chief State Sanitary Inspector of the corresponding territory for a specific Water supply system based on the sanitary and epidemiological situation in the residential place and the water treatment technology applied.

3.5.1. Conspicuous aquatic organisms and surface film are prohibited to be present in the drinking water.

3.6. Radiological safety of the drinking water shall be defined by its conformity to the standards for radiological safety per parameters provided in Table 5.

Table 5

| Parameters | Units of measurement | Radiological safety parameters |
|-------------------------------|----------------------|--------------------------------|
| Consolidated parameters 1) | | |
| Specific total alpha-activity | Bq/kg | 0.2 |
| Specific total beta-activity | Bq/kg | 1.0 |
| Radionuclides 2) | | |
| ²²² Rn (Rn) 3) | Bq/kg | 60 |
| SUM of radionuclides 3) | units | <= 1.0 |

Notes.

1) Should the parameters be exceeded, the test for radionuclides content in the water is conducted.

2) The list of the radionuclides determined in water is established in accordance with the sanitary legislation. Radon determination for the underground water supply sources is obligatory.

3) If several radionuclides are jointly physically present in water the condition $\sum \frac{A_i}{IL_i} \leq 1$ shall be satisfied, where A_i is specific activity of the i -th radionuclide in the water; and IL_i is the corresponding intervention level according to Supplement 2a to SanPiN 2.6.1.2523-09 <*> "Radiological safety standards (NRB-99/2009)". Should the condition be not satisfied, the water assessment is conducted in accordance with the sanitary legislation.

<*> Registered with the Ministry of Justice of the Russian Federation on 14.08.2009, registration number 14534.

(Cl. 3.6 in the version of Amendments No.2 approved by Resolution of Chief State Sanitary Inspector of the Russian Federation No. 10 dated February 25, 2010)

3.6.1. Excluded from May 1, 2010. — Amendments No.2 approved by Resolution of Chief State Sanitary Inspector of the Russian Federation No. 10 dated February 25, 2010.

4. Control Over Drinking Water Quality

4.1. In accordance with the Federal Law "On Sanitary and Epidemiological Well-being of Population" state sanitary and epidemiological supervision and production control shall be executed over the drinking water quality.

4.2. Production control over the drinking water quality shall be conducted by a private entrepreneur or a legal entity operating the water supply system in line with the Work Program.

The private entrepreneur or the legal entity operating the water supply system in line with the Work Program shall constantly control the quality of water at withdrawal points, before the water supply to the distribution network as well as at the water supply points of the external and internal water supply network.

4.3. The amount and frequency of water sampling at withdrawal points collected for laboratory tests are set in compliance with the requirements specified in Table 6.

Table 6

| Parameter types | Amount of samples per one year, at least | |
|----------------------------------|--|---------------------------|
| | For underground sources | For surface sources |
| Microbiological | 4 (as per seasons) | 12 (on the monthly basis) |
| Parasitological | not conducted | - "—" |
| Organoleptic | 4 (as per seasons) | 12 (on the monthly basis) |
| Composite indices | - "—" | - "—" |
| Inorganic and organic substances | 1 | 4 (as per seasons) |
| Radiological | 1 | 1 |

4.4. Types of the tested parameters and the amount of the analyzed samples of drinking water before its supply into the distribution network are set in compliance with the requirements specified in Table 7.

Table 7

| Parameter types | Amount of samples per one year, at least | | | | |
|--|--|---------|----------|---------------------|----------|
| | For underground sources | | | For surface sources | |
| | Size of population provided with water from this Water supply system, thousand people. | | | | |
| | up to 20 | 20—100 | Over 100 | up to 100 | Over 100 |
| Microbiological | 50 (1) | 150 (2) | 365 (3) | 365 (3) | 365 (3) |
| Parasitologic | not conducted | | | 12 (4) | 12 (4) |
| Organoleptic | 50 (1) | 150 (2) | 365 (3) | 365 (3) | 365 (3) |
| Composite indices | 4 (4) | 6 (5) | 12 (6) | 12 (6) | 24 (7) |
| Inorganic and organic substances | 1 | 1 | 1 | 4 (4) | 12 (6) |
| Parameters related to water treatment technology | Residual chlorine, residual ozone—at least once per hour, other reagents - at least once per a shift | | | | |
| Radiological | 1 | 1 | 1 | 1 | 1 |

Notes.

1. The following frequency of the water samples collection is established:

(1)—weekly, (2)—three times per week, (3)—daily, (4)—once per a season, (5)—once per two months, (6)—monthly, (7)—twice per month.

2. Should there be no water sterilization at the underground sources water supply system supplying water to population up to 20 thousand people, the collection of samples for microbiological and organoleptic tests shall be conducted at least once per month.

3. During the flood and emergency situations periods the robust regime for drinking water quality control shall be set subject to the agreement with the State Sanitary and Epidemiological Supervision Center.

4.5. Production control over the quality of drinking water in the distribution water supply network is conducted taking into account with microbiological and organoleptic parameters with the frequency specified in Table 8.

Table 8

| Size of the serviced population, thousand people | Amount of samples per month |
|--|--|
| up to 10 | 2 |
| 10—20 | 10 |
| 20—50 | 30 |
| 50—100 | 100 |
| over 100 | 100 + 1 sample for each 5 thousand people, population of over 100 thousand |

Note.

The samples do not include obligatory control samples collected after the repair and other technical works within the distribution network.

4.6. Collection of samples in the distribution network is conducted from the outdoor water stations at the most elevated and the dead-end parts of the network as well as from the taps of internal water supply networks of all the houses equipped with pumping and local water storage tanks.

4.7. Production control over the drinking water quality in line with the Work Program shall be conducted by laboratories of the private entrepreneurs or legal entities operating the water supply system or subject to agreements with the said persons by laboratories of other organizations accredited within the specified procedure for the right to conduct tests (analyses) of drinking water quality.

4.8. State sanitary and epidemiological supervision over the drinking water quality shall be executed by the bodies and organizations of the State Sanitary and Epidemiological Service in line with the regulatory and methodological documents of the State Sanitary and Epidemiological Service of the Russian Federation within the scheduled procedures and on sanitary and epidemiological indications.

4.9. In order to conduct laboratory tests (analyses) of drinking water quality the metrologically certified standard techniques adopted by the State Committee of the Russian Federation for Standardization and Metrology or the Ministry of Health of the Russian Federation. The collection of samples for analysis shall be conducted in accordance with the requirements of the national standards.

Supplement 1

(obligatory)

RULES FOR

ESTABLISHMENT OF THE CONTROLLED PARAMETERS FOR THE DRINKING WATER QUALITY AND DEVELOPMENT OF THE WORK PROGRAM FOR THE DRINKING WATER QUALITY PRODUCTION CONTROL

I. Procedure for organization of works on selection of the chemical composition parameters for drinking water

1. In accordance with Cl. 3.3 of these Sanitary Rules the selection of chemical composition parameters for drinking water, which are subject to constant production control, shall be conducted for each Water supply system on the basis of the results of the chemical composition of water from water supply sources assessment as well as drinking water technology in the water supply system.

2. Selection of parameters characterizing chemical composition of drinking water for conduction of advanced tests shall be conducted by the organization operating the water supply system together with the State Sanitary and Epidemiological Supervision Center in the city or district within two stages.

2.1. At the first stage the organization operating the water supply system together with the State Sanitary and Epidemiological Supervision Center analyze the following materials for the period of at least three last years:

- national statistic reports of companies and organizations as well as other official data on the composition and volume of the waste water coming into the water supply services above the withdrawal point within their bailing territory;

- materials of the environment protection bodies, weather service, water resources management, geology and subsurface use department, companies and organizations on the quality of the surface, underground and drinking water in the water supply system following the results of the conducted monitoring of water quality and production control;

- materials of the State Sanitary and Epidemiological Supervision Center following the results of the sanitary inspections of companies and organizations conducting business activities and being a source for pollution of the surface and underground water as well as following the results of the tests for water quality in the water use populated places and within the water supply system;

- materials of the bodies for agriculture management and development as to the as the range and gross volume of pesticides and agricultural chemicals used in the territory of water collection (for the surface water source) and within the sanitary protection zone (for the underground water source). Based on the analysis conducted the list of substances characterizing the chemical composition of water from a specific water supply source as well as of hygienic standards in accordance with Supplement 2 of these Sanitary Rules.

2.2. At the second stage the private entrepreneurs and legal entities operating the water supply systems shall conduct advanced laboratory tests of water in line with the prepared list of chemical substances as well as in accordance with the parameters provided in Table 2 of these Sanitary Rules.

2.2.1. At conduction of the advanced tests before water supply into the distribution network for the water supply system with chemical reagent water treating methods of treatment the parameters provided in Table 3 of these Sanitary Rules are additionally included.

2.2.2. Advanced laboratory tests of water are conducted during a year at withdrawal points of the water supply system; and should the system involve water treatment or mix the water from different withdrawal points, the tests are also conducted before drinking water supply to the distribution network.

2.2.3. The minimal amount of the analyzed water samples depending on the type of the water supply source, which provides for the regularity of the water quality receipt within the year, shall be the following:

- for underground sources—4 samples per year, collected once per season;

- for surface sources—12 samples per year collected on the monthly basis.

2.2.4. If it is necessary to obtain more representative and reliable information about the chemical composition of water and changes of the substances concentration therein, the amount of the tested samples and frequency of collection thereof shall be increased in line with the set objectives for the water supply source water quality assessment.

2.2.5. At conduction of the advanced tests it is recommended to use modern general-purpose physical and chemical methods of water media analysis (chromatography-mass spectrometer, etc.) which provide for maximally full information of the chemical composition of water.

2.3. The State Sanitary and Epidemiological Supervision Center shall analyze the results of the advanced tests of the water chemical composition for each Water supply system, and then shall determine the potential hazard of the waterborne chemical substances for the population's health taking into consideration the sanitary and hygienic conditions for the drinking water use by the population and sanitary and epidemiological situation in the city, residential place or district.

2.4. Based on the conducted assessment the State Sanitary and Epidemiological Supervision Center shall develop its suggestions on the list of the controlled parameters, drinking water samples amount and frequency of collection thereof for continuous production control.

II. Procedure for Development of Work Program for Production Control Over Drinking Water Quality

1. Private entrepreneurs and legal entities operating the water supply systems shall develop the Work Program on the basis of these Sanitary Rules.

2. In case of a Water supply system with several withdrawal points the Work Program shall be developed for each withdrawal point taking into account the peculiarities thereof. In case of underground withdrawal points joined in one sanitary protection zone and operating the same water-bearing formation one Work Program can be developed subject to availability of hydrogeological supporting rationale.

3. The Work Program shall contain:

3.1. The list of the controlled parameters of the water quality and the related hygienic standards specified by these Sanitary Rules.

- microbiological and parasitological (Cl. 3.3, Table 1);

- organoleptic (Cl. 3.5, Table 4);

- radiological (Cl. 3.6, Table 5);

- composite (Cl. 3.4.1, Table 2);

- residual amounts of reagents (Cl. 3.4.1, Table 3);

- chemical substances selected for continuous control in accordance with the rules specified in Section 1 of this Supplement (Cl. 3.4.1, Table 2 and Cl. 3.4.3, Supplement 2 to the Sanitary Rules).

3.2. Methods for identification of the controlled parameters.

3.3. Layout of the water samples collection points at the withdrawal points before the water supply into the distribution network of the water system (in the clean-water reservoir) as well as at water supply points of the external and internal water supply network.

3.4. The amount of the controlled water samples and frequency of collection thereof for laboratory analyses (tests), the list of parameters identified in the tested water samples.

3.5. Calendar schedules for water samples collection and conduction of analyses (tests) thereof.

3.6. The amount of the controlled water samples and frequency of collection thereof shall be independently specified for each Water supply system considering the suggestions of the State Sanitary and Epidemiological Supervision Center, though these shall not be less than established in Cl. 4.3, Table 6, Cl. 4.4, Table 7 and Cl. 4.5 Table 8 of these Sanitary Rules.

4. Work Program shall provide for conduction of a monthly analysis of the water quality control results as well as shall specify the procedure for transfer of information about the control results to the water supply system management body, the State Sanitary and Epidemiological Supervision Center and local authorities.

5. Work Program shall be transferred for agreement to the State Sanitary and Epidemiological Supervision Center in the city or district for its further approval within the established procedure.

6. Work Program shall be approved for the period up to 5 years. Subject to agreement with the State Sanitary and Epidemiological Supervision Center modifications and amendments can be introduced into the Work Program within the mentioned period.

(obligatory)

HYGIENIC STANDARDS FOR CONTENT OF HARMFUL SUBSTANCES IN DRINKING WATER

1. This list includes the hygienic standards for harmful substances in the drinking water. The list integrates separate chemical substances which can be present in the drinking water in their specified form and can be identified by modern analytic methods.

2. Chemical substances are arranged in the list in accordance with the structure of organic and inorganic compounds. Each subsection is the extension of the corresponding section. The substances within the subsections are arranged in ascending order of the numerical values of their standards.

If the structure of the organic substance molecule provides for its simultaneous assigning to several chemical class, then the substance is put in the list in accordance with its functional group with the maximum expansion index (as to the horizontal classification).

Organic acids including the pesticides are standardized as to the anion regardless of the form in which the acid is provided in the list (as an acid, acid ion or acid's salt).

Elements and cations (Cl. 1 of "Inorganic substances" section) are standardized cumulatively for all oxidation rates unless provided otherwise.

3. The list has the following vertical classification:

3.1. The first column of the list indicates the most frequently used names of chemical substances.

3.2. The second column provides synonyms of the chemical substances names and some common and generally accepted names.

3.3. The third column shows the MAC and the approximate permissible level (APL) in mg/l, where:

MAC is the maximum allowed concentration at which the substances have no direct or indirect influence on human health condition (subject to influence on the human organism for the whole life) and do not impair the hygienic conditions of water use;

APL (marked by the asterisk) are the approximate permissible levels of substances in the tap water determined on the basis of computational, express and experimental methods for the toxicity level forecast.

If the column of the standard level indicates «"None"», it means that the concentration of such a compound in drinking water shall be less than the detection level of the testing method applied.

3.4. The fourth column specifies the limiting characteristic for the substances' hazard in accordance with which the standard was set:

- "s.-t."—sanitary and toxicological;

- "org"—organoleptic with explanation of the water organoleptic attributes behavior ("odr"—changes the water odor; "clr"—colors the water; "fmg"—provides for foaming; "fm"—makes a film on the surface of the water; "flavr"—gives the water a flavour; "op"—produces opalescence).

3.5. The fifth column demonstrates the substance hazard category:

1st category—extremely hazardous;

2nd category—highly hazardous;

3rd category—hazardous;

4th category—moderately hazardous.

The classification is based on the parameters characterizing the different hazard level of chemical substances contaminating drinking water for people depending on toxicity, cumulativeness, ability to give rise to delayed actions and the hazard limiting value (HLV).

Substances hazard categories shall be considered at:

- selection of compounds to be primarily controlled in the drinking water;

- establishment of sequence of the water-protective measures requiring additional financial investments;

- substantiation of recommendations on replacement of highly hazardous substances to less hazardous in production processes;

- determining the priority of development of the selective methods for analytic control of substances in water.

HYGIENIC STANDARDS FOR

CONTENT OF HARMFUL SUBSTANCES IN DRINKING WATER

| Substance name | Synonyms | Standard value in mg/l | Hazard parameter | Hazard category |
|-----------------------------|----------|------------------------|------------------|-----------------|
| 1 | 2 | 3 | 4 | 5 |
| Inorganic Substances | | | | |
| 1. Elements, cations | | | | |
| Thallium | | 0.0001 | s.-t. | 2 |
| Phosphorus, elemental | | 0.0001 | s.-t. | 1 |
| Niobium | | 0.01 | s.-t. | 2 |
| Tellurium | | 0.01 | s.-t. | 2 |
| Samarium | | 0.024 <1> | s.-t. | 2 |
| Lithium | | 0.03 | s.-t. | 2 |
| Stibium | | 0.05 | s.-t. | 2 |
| Wolframium | | 0.05 | s.-t. | 2 |
| Argentum | | 0.05 | s.-t. | 2 |
| Vanadium | | 0.1 | s.-t. | 3 |
| Bismuth | | 0.1 | s.-t. | 2 |
| Cobalt | | 0.1 | s.-t. | 2 |
| Rubidium | | 0.1 | s.-t. | 2 |
| Europium | | 0.3 <1> | org. flavr. | 4 |
| Ammonia (as to nitro) | | 2.0 | s.-t. | 3 |
| Chromium(Cr ³⁺) | | 0.05 | s.-t. | 3 |
| Silicium | | 10.0 | s.-t. | 2 |
| Sodium | | 200.0 | s.-t. | 2 |
| 2. Anions | | | | |
| Rhodanide ion | | 0.1 | s.-t. | 2 |
| Chlorite ion | | 0.2 | s.-t. | 3 |

| | | | | |
|------------------------|------------------------|-------|----------------|---|
| Bromide ion | | 0.2 | s.-t. | 2 |
| Persulphate ion | | 0.5 | s.-t. | 2 |
| Hexanitrocobaltate ion | | 1.0 | s.-t. | 2 |
| Ferrocyanide ion | | 1.25 | s.-t. | 2 |
| Hydrosulphide ion | | 3.0 | s.-t. | 2 |
| Nitrite ion | | 3.0 | org. | 2 |
| Perchlorate ion | | 5.0 | s.-t. | 2 |
| Chlorate ion | | 20.0 | org. flavr. | 3 |
| Sulphurated hydrogen | Hydrogen sulphide | 0.003 | org. odr. | 4 |
| Hydrogen dioxide | Hydrogen peroxide | 0.1 | s.-t. | 2 |
| Organic Substances | | | | |
| 1. Carbohydrates | | | | |
| 1.1. aliphatic | | | | |
| Isoprene | 2-methyl butal, 3diene | 0.005 | org. odr. | 4 |
| Butadiene-1,3 | Divinyl | 0.05 | org. odr. | 4 |
| Butylene | But-1-ene | 0.2 | org. odr. | 3 |
| Ethylene | Ethene | 0.5 | org. odr. | 3 |
| Propylene | Propene | 0.5 | org. odr. | 3 |
| Isobutene | 2-methylprop-1-ene | 0.5 | org. odr. | 3 |
| 1.2. Closed-chain | | | | |
| 1.2.1. Alicyclic | | | | |
| 1.2.1.1. Mononucleate | | | | |
| Cyclohexene | tetrohydrobenzene | 0.02 | s.-t. | 2 |
| Cyclohexane | Hexahydrobenzene, he- | 0.1 | s.-t. | 2 |

| | | | | |
|------------------------|---|-------|-------------|---|
| | xamethylene | | | |
| 1.2.1.2. Multinucleate | | | | |
| Norbornene | 2,3-Dicyclo (2.2.1)heptene | 0.004 | org. odr. | 4 |
| Dicycloheptadiene | Bicyclo (2,2,1)hepta-2,5-diene, norbornadiene | 0.004 | org. odr. | 4 |
| Dicyclopentadiene | Tricyclopentadiene-3,8-diene 3a,4,7,7a-tetrahydro-4,7-methano-1 H-inden | 0.015 | org. odr. | 3 |
| 1.2.2. Aromatic | | | | |
| 1.2.2.1. Mononucleate | | | | |
| Benzene | | 0.01 | s.-t. | 2 |
| Ethylbenzene | | 0.01 | org. flavr. | 4 |
| m-Diethylbenzene | 1,3-Diethylbenzene | 0.04 | org. odr. | 4 |
| Xylene | Dimethylbenzene | 0.05 | org. odr. | 3 |
| Diisopropylbenzene | di-1-methylethylbenzene | 0.05 | s.-t. | 2 |
| Monobenzyltoluene | 3-Benzyltoluene | 0.08 | org. odr. | 2 |
| Butylbenzene | 1-Phenylbutane | 0.1 | org. odr. | 3 |
| Isopropylbenzene | cumene, 1-methylethylbenzene | 0.1 | org. odr. | 3 |
| Styrole | Vinylbenzene | 0.1 | org. odr. | 3 |
| alpha-Methylstyrene | (1-Methylvinyl)benzene | 0.1 | org. flavr. | 3 |
| Propyl benzene | 1-Phenypropane | 0.2 | org. odr. | 3 |

| | | | | |
|--------------------------------------|---|-----------|-----------|---|
| n-t-butyltoluene | 1-(1,1-Dimethylethyl) 4-methylbenzene, 1 1-methyl-4-t-butylbenzen | 0.5 | org. odr. | 3 |
| Toluene | Methylbenzene | 0.5 | org. odr. | 4 |
| Dibenzyltoluene | [(3-Methyl-4-benzyl)phenyl]phenylmethane | 0.6 | org. odr. | 3 |
| 1.2.2.2. Multinucleate | | | | |
| Benz(a)pyrene | | 0.000-005 | s.-t. | 1 |
| 1.2.2.2.1. biphenyls | | | | |
| Diphenyl | Biphenyl, phenylbenzene | 0.001 | s.-t. | 2 |
| Alkylenediphenyl | | 0.4 | org. film | 2 |
| 1.2.2.2.2. condensated | | | | |
| Naphthalene | | 0.01 | org. odr. | 4 |
| 2. Halogen-containing compounds | | | | |
| 2.1. Aliphatic | | | | |
| 2.1.1. Containing only maximum bonds | | | | |
| Iodoform | Triiodomethane | 0.0002 | org. odr. | 4 |
| Tetrachloroheptane | | 0.0025 | org. odr. | 4 |
| 1,1,1,9-Tetrachlorononane | | 0.003 | org. odr. | 4 |
| Butyl chloride | 1-chlorbutane | 0.004 | s.-t. | 2 |
| 1,1,1,5-Tetrachloropentane | | 0.005 | org. odr. | 4 |
| Tetrachloride carbon | Tetrachloromethane | 0.006 | s.-t. | 2 |

| | | | | |
|------------------------------------|-------------------------------------|-------|-----------|---|
| 1,1,1,11-Tetrachlorohendecane | | 0.007 | org. odr. | 4 |
| Hexachlorobutane | | 0.01 | org. odr. | 3 |
| Hexachloroethane | | 0.01 | org. odr. | 4 |
| 1,1,1,3-Tetrachloropropane | | 0.01 | org. odr. | 4 |
| 1-Chloro-2,3-dibromopropane | 1,2-Dibrom-3-chloropropane, nemagon | 0.01 | org. odr. | 3 |
| 1,2,3,4-Tetrachlorobutane | | 0.02 | s.-t. | 2 |
| Pentachlorobutane | | 0.02 | org. odr. | 3 |
| Perchlorobutane | | 0.02 | org. odr. | 3 |
| Pentachloropropane | | 0.03 | org. odr. | 3 |
| Dichlorobromomethane | | 0.03 | s.-t. | 2 |
| Chlorodibromomethane | | 0.03 | s.-t. | 2 |
| 1,2-Dibromo-1,1,5-trichloropentane | Ethyl bromide | 0.04 | org. odr. | 3 |
| 1,2,3-Trichloropropane | | 0.07 | org. odr. | 3 |
| Trifluorochloropropane | Freon 253 | 0.1 | s.-t. | 2 |
| 1,2-Dibromopropane | | 0.1 | s.-t. | 3 |
| Bromoform | Tribromomethane | 0.1 | s.-t. | 2 |
| Tetrachloroethane | | 0.2 | org. odr. | 4 |
| Chloroethyl | Chloroethane ethyl-chloride, | 0.2 | s.-t. | 4 |
| 1,2-Dichloropropane | | 0.4 | s.-t. | 2 |
| 1,2-Dichloroisobutane | 2-Methyl-1,2 dichloropropane | 0.4 | s.-t. | 2 |
| Dichloromethane | Methylene chloride | 7.5 | org. odr. | 3 |

| | | | | |
|-------------------------------------|---------------------------------|----------|-------------|---|
| Difluorochloromethane | Freon-22 | 10.0 | s.-t. | 2 |
| Difluorodichloromethane | Freon-12 | 10.0 | s.-t. | 2 |
| Methyl chloroform | 1,1,1-tri-chloroethane | 10.0 <1> | s.-t. | 2 |
| 2.1.2. Containing double bonds | | | | |
| Tetrachloropropene | | 0.002 | s.-t. | 2 |
| 2-Methyl-3-chloroprop-1-ene | Methallyl chloride | 0.01 | s.-t. | 2 |
| beta-Chloroprene | 2-chlorobuta-1,3-diene | 0.01 | s.-t. | 2 |
| Hexachlorobutadiene | Perchlorobuta-1,3-diene | 0.01 | org. odr. | 3 |
| 2,3,4-Trichlorobutene-1 | 2,3,4-Tri-chlorobut-1-ene | 0.02 | s.-t. | 2 |
| 2,3-Dichlorobutadiene-1,3 | 2,3-Dichlorobuta-1,3-diene | 0.03 | s.-t. | 2 |
| 1,1,5-Trichloropentene | | 0.04 | org. odr. | 3 |
| Vinyl chloride | Chloroethene chloroethylene | 0.05 | s.-t. | 2 |
| 1,3-Dichlorobutene-2 | 1,3-Dichlorobut-2-ene | 0.05 | org. odr. | 4 |
| 3,4-Dichlorobutene-1 | | 0.2 | s.-t. | 2 |
| Allyl chloride | 3-chloroprop-1-ene | 0.3 | s.-t. | 3 |
| 1,1-Dichloro-4-methylpentadiene-1,4 | Diene-1,4 | 0.37 | org. flavr. | 3 |
| Dichloropropene | | 0.4 | s.-t. | 2 |
| 3.3-Dichloroisobutylene | 3.3-Dichloro-2-methyl-1-propene | 0.4 | s.-t. | 2 |
| 1,3-Dichloroisobutylene | 2-Methyl-1,3-dichloroprop | 0.4 | s.-t. | 2 |

| | | | | |
|---|---|-------|-------------|---|
| | -1-ene | | | |
| 1,1-Dichloro-4-methylpentadiene-1.3 | Diene-1,3 | 0.41 | org. odr. | 3 |
| 2.2. Closed-chain | | | | |
| 2.2.1. Alicyclic | | | | |
| 2.2.1.1. Mononucleate | | | | |
| Hexachlorocyclopentadiene | 1,2,3,4,5,5-Hexachloro-1,3-cyclopentadiene | 0.001 | org. odr. | 3 |
| 1,1-Dichlorocyclohexane | | 0.02 | org. odr. | 3 |
| 1,2,3,4,5,6-Hexachlorocyclohexane | Hexachlorane | 0.02 | org. odr. | 4 |
| Perchloromethylene-cyclopentene | 4-(Dichloromethylene)-1,2,3,3,5,5-Hexachlorocyclopentene | 0.05 | org. odr. | 4 |
| Chlorocyclohexane | | 0.05 | org. odr. | 3 |
| 2.2.1.2. Multinucleate | | | | |
| 1,2,3,4,10,10-Hexachloro-1,4,4a,5,8,8a-hexahydro-1,4-endo-exo-5,8-dimethanenaphthalene | 1,4,4a,5,8,8a-Hexahydro-1,2,3,4,10,10-hexachloro-1,4,5,8-dimethanenaphthalene, aldrin | 0.002 | org. flavr. | 3 |
| 1,4,5,6,7,8,8-Heptachlor-4,7-endo-methylene-3a,4,7,7a-tetrahydroindene-tetrahydroindene | 3a,4,7,7a-Tetrahydro-1,4,5,6,7,8,8-heptachloro-4,7-methano-1H-indene, heptachlor | 0.05 | s.-t. | 2 |
| beta-Dihydroheptachlor | 2,3,3a,4,7,7a-Hexahydro-2,4,5,6,7,8,8-heptachlor- | 0.1 | org. odr. | 4 |

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|--|--|-------|-----------|---|
| | 4,7-methanoindene, dichlor | | | |
| Polychloropinene | | 0.2 | s.-t. | 3 |
| 2.2.2. Aromatic | | | | |
| 2.2.2.1. Mononucleate | | | | |
| 2.2.2.1.1. with halogen atom in the nuclear | | | | |
| 2,5-Dichloro-n-t-butyltoluene | 1,4-Dichloro-2-(1,1-dimethyl)-5-methylbenzene | 0.003 | org. odr. | 3 |
| o-Dichlorobenzene | 1,2-Dichlorobenzene | 0.002 | org. odr. | 3 |
| Chlor-n-t-butyltoluene | 1-Methyl-4-(1,1-dimethylethyl)-2benzene chloride | 0.002 | org. odr. | 4 |
| 1,2,3,4-tetrachlorobenzene | | 0.01 | s.-t. | 2 |
| Chlorobenzene | | 0.02 | s.-t. | 3 |
| 2,4-Dichlorotoluene | 2,4-Dichlor-1-methylbenzene | 0.03 | org. odr. | 3 |
| 1,3,5-Trichlorobenzene | | 0.03 | org. odr. | 3 |
| 2,3,6-Trichlorotoluene | | 0.03 | org. odr. | 3 |
| o- and n-Chlorotoluene | o- and n-chloro-methylbenzene | 0.2 | s.-t. | 3 |
| 2,3,6-Trichlor-n-t-butyltoluene | | 0.1 | org. odr. | 4 |
| 2.2.2.1.2. with halogen atom in the side chain | | | | |
| Benzyl chloride | Chloromethylbenzene | 0.001 | s.-t. | 2 |
| Hexachlorometaxylene | 1,3-Bis(trichloro- | 0.008 | org. odr. | 4 |

| | | | | |
|---------------------------------|--|-------|-----------|---|
| | methyl) benzene | | | |
| Hexachloroparaxylene | 1.4-Bis(tri-chloro-methyl) benzene | 0.03 | org. odr. | 4 |
| Benzotrichloride | Trifluoro-methyl-benzene | 0.1 | s.-t. | 2 |
| 2.2.2.2. Multinucleate | | | | |
| 2.2.2.2.1. Biphenyls | | | | |
| Monochlorodiphenyl | Monochloro-biphenyl | 0.001 | s.-t. | 2 |
| Dichlorodiphenyl | Dichlorobi-phenyl | 0.001 | s.-t. | 2 |
| Trichlorodiphenyl | Trichlorobi-phenyl | 0.001 | s.-t. | 1 |
| Pentachlorodiphenyl | Pentachloro-biphenyl | 0.001 | s.-t. | 1 |
| 2.2.2.2.2. condensated | | | | |
| 2-2-Chloronaphthalene | | 0.01 | org. odr. | 4 |
| 3. Oxygen containing compounds | | | | |
| 3.1. Alcohols and simple ethers | | | | |
| 3.1.1. Monohydroxy alcohols | | | | |
| 3.1.1.1. Aliphatic alcohols | | | | |
| 3-Methyl-3-butenelol | Isobutenyl-carbinol | 0.004 | s.-t. | 2 |
| Heptyl alcohol, regular | Heptan-1-ol hexyl-carbinol | 0.005 | s.-t. | 2 |
| 3-Metal-1-butene-3-ol | 2-Methylpro-2-pen-1-ol, dimethylvi-nylcarbinol, isoprene alcohol | 0.005 | s.-t. | 2 |
| Hexyl alcohol, regular | Hexan-1-ol, amylcarbi- | 0.01 | s.-t. | 2 |

| | | | | |
|--------------------------|---|------|-------------|---|
| | nol, pentyl-carbinol | | | |
| Hexyl alcohol, secondary | 1-Methylpentan-1-ol, hexan-2-ol, methylbuten-carbinol | 0.01 | s.-t. | 2 |
| Hexyl alcohol, tertiary | 2-Methylpentan-2-ol, diethylmethylcarbinol flotation reagent, TTS | 0.01 | s.-t. | 2 |
| Nonyl alcohol, regular | Nonan-1-ol, octylocarbinol | 0.01 | s.-t. | 2 |
| Octyl alcohol , regular | Octan-1-ol, heptylcarbinol | 0.05 | org. flavr. | 3 |
| Butyl alcohol, regular | Butan-1-ol, propylcarbinol | 0.1 | s.-t. | 2 |
| Allyl alcohol | Prop-2-en-1-ol, vinyl-carbinol | 0.1 | org. flavr. | 3 |
| Isobutyl alcohol | 2-Methylpropan-1-ol, isopropyl-carbinol | 0.15 | s.-t. | 2 |
| Butyl alcohol, secondary | Butan-2-ol, methylisobutyl-carbinol | 0.2 | s.-t. | 2 |
| Propyl alcohol | Propan-1-ol, ethyl-carbinol | 0.25 | org. odr. | 4 |
| Isopropyl alcohol | Propan-2-ol, dimethylcarbinol | 0.25 | org. odr. | 4 |
| Butyl alcohol, tertiary | t-butyl alcohol , 1,1-dimethyl ethanol, trimethylcarbinol, 2-me- | 1.0 | s.-t. | 2 |

| | | | | |
|---|---|------|-----------|---|
| | thylpropan-2-ol | | | |
| Amyl alcohol | Pentan-1-ol, butylkcarbinol | 1.5 | org. odr. | 3 |
| Methyl alcohol | Methanol, carbinol | 3.0 | s.-t. | 2 |
| 3.1.1.1.1. Halogen substituted monohydroxy alcohols | | | | |
| Ethylene chlorhydrin | 1-Chlor-2-hydroxyethane, 2-chloroethanol, 2-chloroethyl alcohol, chloromethylcarbinol, 1-chloroethan-2-ol | 0.1 | s.-t. | 2 |
| 1,1,7-Trihydro dodecafluoroheptyl alcohol | P-3 | 0.1 | org. odr. | 4 |
| 1,1,3- Trihydro tetradecafluoropropyl alcohol | P-1 | 0.25 | org. odr. | 3 |
| 1,1,5- Trihydro octafluoropentylalcohol | P-2 | 0.25 | org. odr. | 4 |
| 1,1,9- Trihydro hexadecafluorononyl alcohol | P-4 | 0.25 | org. odr. | 4 |
| 1,1,13- Trihydro tetraeicosafluorotridecyl alcohol | P-6 | 0.25 | org. odr. | 3 |
| 1,1,11- Trihydro eicosafluoroundecylic alcohol | P-5 | 0.5 | org. odr. | 3 |
| beta,beta-Dichlorisopropyl alcohol | 1,3-Dichloropropan-2-ol, dichlorohydrin, dichloromethylcarbinol | 1.0 | org. odr. | 3 |
| 1,1-Dihydroperfluoroheptyl alcohol | 2,2,3,3,4,4,5,5,6,6,7,7, | 4.0 | s.-t. | 2 |

| | | | | |
|---|---|-------|-------------|---|
| | 7-Tridecafluoroheptan-1-ol | | | |
| 3.1.1.2. Closed-chain | | | | |
| 3.1.1.2.1. Alicyclic | | | | |
| Cyclohexanol | Hexahydrophenol | 0.5 | s.-t. | 2 |
| 3.1.1.2.2. Aromatic | | | | |
| 3.1.1.2.2.1. Mononucleate | | | | |
| 3.1.1.2.2.1.1. Phenols | | | | |
| Phenol | | 0.001 | org. odr. | 4 |
| m- and n-Cresol | m- and n-Methylphenol, 1-hydroxy-2 (and 4) methylphenol | 0.004 | s.-t. | 2 |
| o- and n-Propylphenol | 1-Hydroxy-2 (and 4)-propyl benzene | 0.01 | org. odr. | 4 |
| Alkylenephenol | | 0.1 | org. | 3 |
| Dimethylphenol | Xylenol | 0.25 | org. odr. | 4 |
| 3.1.1.2.2.1.1.1. halogen substituted | | | | |
| Chlorophenol | | 0.001 | org. odr. | 4 |
| Dichlorophenol | | 0.002 | org. flavr. | 4 |
| Trichlorophenol | | 0.004 | org. flavr. | 4 |
| 3.1.1.2.2.1.2. Containing a hydroxy group in the side chain | | | | |
| 3.1.1.2.2.1.2.1. Halogen substituted | | | | |
| 3.1.1.2.2.2. Condensated | | | | |
| alpha-Naphthol | Napht-1-ol, 1-naphthol | 0.1 | org. odr. | 3 |
| 3-Naphthol | Napht-1-ol, 2-naphthol | 0.4 | s.-t. | 3 |

| | | | | |
|--|--|----------|-------------|---|
| 3.1.2. Simple ethers | | | | |
| 3.1.2.1. Aliphatic | | | | |
| Ethylvinylbutyl-ether | 1-Butoxi-but-1-en-3-ine, butoxi-butenine | 0.002 | org. odr. | 4 |
| Diethylacetal | 1,1-Diethoxoethane | 0.1 | org. odr. | 4 |
| Ethoxylate for pr. alch. C12 - C15 | | 0.1 | org. foam | 4 |
| Diethylether | Ethoxyethane | 0.3 | org. flavr. | 4 |
| Dimethyl ether | Methoximethane | 5.0 | s.-t. | 4 |
| 3.1.2.1.1. Halogen substituted | | | | |
| beta,beta-Dichlordiethyl ether | 1,1'-Oxibis (2-chloroethane), chlorex | 0.03 <1> | s.-t. | 2 |
| 3.1.2.2. Aromatic | | | | |
| Diphenylolpropane | 4,4'-Isopropylidendi-phenol | 0.01 | org. flavr. | 4 |
| m-Phenoxitoluene | 3-Phenoxitoluene | 0.04 | org. | 4 |
| Anisol | Methoxybenzene | 0.05 | s.-t. | 3 |
| 3.1.3. Polyatomic alcohols and mixed compounds | | | | |
| 3.1.3.1. Aliphatic polyatomic alcohols | | | | |
| 2-Methyl2,3-butandiol | Methylbutandiol | 0.04 | s.-t. | 2 |
| Glycerol | Trioxipropane, propantriol | 0.06 <1> | org. foam | 4 |
| Pentaerythritol | 2,2-Dimethylolpropandiol-1,3 | 0.1 | s.-t. | 2 |
| Ethylene glycol | Ethan-1,2- | 1.0 | s.-t. | 3 |

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| | diol | | | |
| 1,4-Butindiol | But-2-in-1,4-diol | 1.0 | s.-t. | 2 |
| 1,4-Butandiol | Butan-1,4-diol | 5.0 | s.-t. | 2 |
| 3.1.3.1.1. Halogen substituted | | | | |
| Monochlorohydrin | 3-Chloropropan-1,2-diol, alpha-chlorohydrine | 0.7 | org. flavr. | 3 |
| 3.1.3.2. Polyatomic phenols | | | | |
| Pyrocatechol | 1,2-Benzenediol, 1,2-dioxibenzene | 0.1 | org. clr. | 4 |
| Pyrogallol | 1,2,3-Trioxibenzene | 0.1 | org. clr. | 3 |
| Hydroquinone | 1,4-Dioxibenzene | 0.2 | org. clr. | 4 |
| 5-Methylresorcine | 5-Methyl,3-benzenediol | 1.0 | org. clr. | 4 |
| 3.1.3.2.1. Halogen substituted | | | | |
| 2,2-Bis-(4-hydroxi-3,5-dichlorphenyl)propane | Tetrachlorodiane | 0.1 | org. flavr. | 4 |
| 3.1.3.3. Containing hydroxy and oxy groups | | | | |
| 3.1.3.3.1. Aliphatic | | | | |
| 2-allyloxiethyl alcohol | | 0.4 | s.-t. | 3 |
| Diethylene glycol | 2,2'-Oxidiethanol | 1.0 | s.-t. | 3 |
| Tetraethylene glycol | 2,2'-Oxidiethylendioxidiethanol | 1.0 | s.-t. | 3 |
| Pentaethylene glycol | 3,6,9,12-Tetraoxatetradecan-1.14-diol, | 1.0 | s.-t. | 3 |

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| | ethylenglycoltetraoxididiethyl ether | | | |
| 3.1.3.3.2. Aromatic | | | | |
| 3-Phenoxibenzyl alcohol | 3-Phenoxiphenylmethanol, 3-phenoxiphenylcarbinol | 1.0 <1> | s.-t. | 3 |
| 3.2. Aldehydes and ketones | | | | |
| 3.2.1. Containing only one oxy group | | | | |
| 3.2.1.1. Aliphatic | | | | |
| 3.2.1.1.1. Aliphatic compounds containing only maximum bonds | | | | |
| Diethylketone | Pentan-3-one 3-oxopentane | 0.1 | org. odr. | 4 |
| Methylethylketone | Butan-2-one, 2-oxobutane | 1.0 | org. odr. | 3 |
| 3.2.1.1.1.1. Halogen substituted | | | | |
| Chloral | Trichloroacetaldehyde | 0.2 | s.-t. | 2 |
| Perfluoroheptanalhydrate | | 0.5 | s.-t. | 2 |
| 3.2.1.1.1.2. Containing hydroxy and oxy groups | | | | |
| Diacetone alcohol | 4-Hydroxi-4-methylpenten-2-one | 0.5 <1> | s.-t. | 2 |
| 3.2.1.1.2. Containing a double bond | | | | |
| Acrolein | Propenal, allyl aldehyde | 0.02 | s.-t. | 1 |
| Mesityl oxide | 2-Methylpent-2-en-4-one | 0.06 <1> | s.-t. | 2 |
| alpha-Ethyl-beta-acrolein | 2-Ethylhexenal | 0.2 | org. odr. | 4 |
| beta-Methylacrolein | But-2-enal, croton | 0.3 | s.-t. | 3 |

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| | aldehyde, 2-butenal | | | |
| 3.2.1.2. Closed-chain | | | | |
| 3.2.1.2.1. Alicyclic | | | | |
| Cyclohexanol | | 0.2 | s.-t. | 2 |
| 3.2.1.2.1.1. Halogen substituted | | | | |
| Bromcamphora | | 0.5 <1> | org. odr. | 3 |
| 3.2.1.2.2. Aromatic | | | | |
| 3.2.1.2.2.1. Containing mononucleate aromatic substitutes | | | | |
| m-Phenoxibenzaldehyde | 3-Phenoxy- benzaldehyde | 0.02 | s.-t. | 2 |
| Acetophenone | | 0.1 | s.-t. | 3 |
| 2,2-Dimethoxy-1,2-diphenylethanone | 2,2-Dimethoxy-2-phenylacetophenone | 0.5 <1> | org. odr. | 3 |
| 3.2.1.2.2.1.1. Halogen substituted | | | | |
| m-Brombenzaldehyde | 3-Brombenzaldehyde | 0.02 | s.-t. | 2 |
| Pentachloracetophenone | 1-(Pentachlorophenyl)ethanone | 0.02 | org. flavr. | 3 |
| 3,3-Dimethyl-1-chlor-1-(4-chlorphenoxi)butan-2-one | | 0.04 | s.-t. | 4 |
| 3.2.2. Containing more than one oxy group | | | | |
| Tetrahydroquinone | Cyclohexan-1,4-dione, 1,4-dioxocyclohexane | 0.05 | org. odr. | 3 |
| Glutaric aldehyde | Glutaric dialdehyde | 0.07 | s.-t. | 2 |
| Acetylacetonates | | 2.0 <1> | s.-t. | 2 |
| Anthraquinones | 9,10-Dihydro-9,10-dioxoanthracene, 9,10-anthracendione | 10,0 | s.-t. | 3 |

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| 3.2.2.1. Halogen substituted | | | | |
| 2,3,5,6-Tetrachlor-n-benzoquinone | Chloranil, tetrachlor-quinone | 0.01 | org. clr. | 3 |
| 2,3-Dichlor-5-dichlormethylene-2-cyclopenten-1,4-dione | 4,5-Dichlor-2-(dichlormethylene)-4-cyclopentene-1,3-dione, diketone | 0.1 | org. odr. | 3 |
| 2,3-Dichlor-1,4-naphtoquinone | | 0.25 | s.-t. | 2 |
| 1-Chloranthraquinone | | 3.0 | s.-t. | 2 |
| 2-Chloranthraquinone | beta-Chloranthraquinone | 4.0 | s.-t. | 2 |
| 3.2.2.2. containing a hydroxy group | | | | |
| 1,5-Dihydroxyanthraquinone | 1,5-Dihydroxi-9,10-anthracendione | 0,1 | org. clr. | 3 |
| 1,8-Dihydroxianthraquinone | Dantrone | 0.25 | org. clr. | 3 |
| 1,2-Dihydroxianthraquinone | 1,2-Dihydroxi-9,10-anthracendione, alizarin | 3.0 | s.-t. | 2 |
| 1,4,5,8-Tetrahydroxianthraquinone | 1,4,5,8-Tetrahydroxi-9,10-anthracendione | 3.0 | s.-t. | 2 |
| 1,4-Dihydroxianthraquinone | Quinizarine | 4.0 | s.-t. | 2 |
| 3.3. Carboxylic acids and their derivative compounds | | | | |
| 3.3.1. Carboxylic acids and their ions | | | | |
| 3.3.1.1. Containing one carboxy group | | | | |
| 3.3.1.1.1. Aliphatic | | | | |
| 3.3.1.1.1.1. Containing only maximum bonds | | | | |

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| Stearinic acid, salt | octadecanoic acid, salt | 0.25 <1> | org. susp. | 4 |
| 3.3.1.1.1.1.1. Halogen substituted | | | | |
| alpha, alpha, beta-trichloropropionic acid | 2,2,3-Tri-trichloropropionic acid | 0.01 | org. flavr. | 4 |
| Chlorenantic acid | 7-chloro-heptanic acid | 0.05 | org. odr. | 4 |
| Monochloroacetic acid, salt | Chloroacetic acid, salt | 0.05 | s.-t. | 2 |
| Chloroundecanoic acid | 11-undecanoic acid | 0.1 | org. odr. | 4 |
| Chlorpelargonic acid | 9-chlor-nonanic acid | 0.3 | org. odr. | 4 |
| Perfluorovaleric acid | Nonafluoropentanic acid, perfluoropentanic acid | 0.7 | s.-t. | 2 |
| alpha-Monochloropropionic acid | 2-Chlorpropionic acid | 0.8 | org. flavr. | 3 |
| Hydroperfluoroenantic acid | 2,2,3,3,4,4,5,5,6,6,7,7-dodecaftoheptanic acid | 1.0 | s.-t. | 2 |
| Perfluoroenantic acid | Perfluoroheptanic acid | 1.0 | s.-t. | 2 |
| 2,2-dichloropropionic acid, sodium salt | Dalapon | 2.0 | org. odr. | 3 |
| Trichloroacetic acid, salt | | 5.0 | org. odr. | 4 |

| 3.3.1.1.1.1.2. Containing aromatic substitutes | | | | |
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| 3.3.1.1.1.1.3. containing hydroxi-, oxi- and oxogroups | | 2.0 | | |
| 5-(2,5-dimethylphenoxi)-2,2-dimethylpentanic acid | Gemfibrozil | 0.001 | s.-t. | 1 |
| Phenoxiacetic acid | Glycolic acid, phenide hydroxi-acetic acid, phenide | 1.0 | s.-t. | 2 |
| 2-(alphanaphtoxi)-propionic acid | 2-(1-naphtalinyloxi)propionic acid | 2.0 | s.-t. | 2 |
| 3.3.1.1.1.1.3.1. Halogen substituted | | | | |
| 2,4-dichlorphenoxi alpha-oleic acid | 4-(2,4-dichlorphenoxi)oleic acid, 2,4DM | 0.01 | s.-t. | 2 |
| 2-Methyl-4-chlorphenoxioleic acid | 4-(2-methylphenoxi)-4-chlorbutanic acid, tropotox | 0.03 | org. odr. | 3 |
| 2,4-Dichlorphenoxi-alpha-propionic acid | 2-(2,4-dichlorphenoxi)propionic acid 2,4-DP | 0.5 | org. flavr. | 3 |
| 3.3.1.1.1.2. Containing olefinic bonds | | | | |
| Acrylic acid | Propan-2-en-carbonic acid | 0.5 | s.-t. | 2 |
| Metacrylic acid | 2-methylpropan-2-en-carbonic acid | 1.0 | s.-t. | 3 |

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| 3.3.1.1.1.2.1. Oxo and halogen containing compounds | | | | |
| alpha, beta-dichlor-beta-forminacrylic acid | 4-oxo-2,3-dichlorisocrotonic acid, mucochloric acid | 1.0 | s.-t. | 2 |
| 3.3.1.1.2. Closed-chain | | | | |
| 3.3.1.1.2.1. Alicyclic | | | | |
| Chrisanthemic acid, salt | 2,2-dimethyl 3-propenyl-1-cyclopropanecarbonic acid, salt; 3-Isobutenyl 2,2-dimethyl 1-cyclopropan-carbonic acid, salt | 0.8 | s.-t. | 3 |
| Naphthenic acids | | 1.0 | org. odr. | 4 |
| 3.3.1.1.2.2. Aromatic | | | | |
| Benzene monocarbonic acid, salt | | 0.6 | org. flavr. | 4 |
| 3.3.1.1.2.2.1. Halogen substituted | | | | |
| 2-chlorbenzene carbonic acid | o-chlorbenzene carbonic acid | 1.0 | org. flavr. | 4 |
| 4-chlorbenzene 4-carbonic acid | n-chlorbenzene carbonic acid | 0.2 | org. flavr. | 4 |
| 2,3,6- trichlorbenzene carbonic acid | | 1.0 | s.-t. | 2 |
| 3.3.1.1.2.2.2. containing hydroxy, oxy and oxo groups | | | | |
| 2-hydroxi-3,6-dichlor benzene carbonic acid | | 0.5 | org. clr. | 3 |
| 2-Metoxy-3,6-dichlor-dichlorbenzene carbonic acid | 2-Metoxy-3,6 dichlorbenzene carbonic acid, | 15.0 | s.-t. | 2 |

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| 3.3.1.2. Polyatomic acids | | | | |
| 3.3.1.2.1. Aliphatic | | | | |
| Maleinic acid | Cisbutendio- ic acid | 1.0 | org. odr. | 4 |
| Adipic acid, salt | Hexandioic acid, salt; 1,4-butan- dicarbonic acid, salt | 1.0 | s.-t. | 3 |
| Sebacylic acid | 1,8-Octan- dicarbonic acid | 1.5 | s.-t. | 3 |
| 3.3.1.2.2. Aromatic | | | | |
| 3.3.1.2.2.1. Halogen substituted | | | | |
| 3.3.2. Compound ethers | | | | |
| 3.3.2.1. Compound ethers of monoatomic acids | | | | |
| 3.3.2.1.1. Aliphatic | | | | |
| 3.3.2.1.1.1. Maximum | | | | |
| 3.3.2.1.1.1.1. Unsubstituted | | | | |
| 3.3.2.1.1.1.1.1. Alcohols containing only maximum bonds | | | | |
| Methyl acetate | Acetic acid, methyl ether; methyl ether of acetic acid | 0.1 | s.-t. | 3 |
| Ethyl acetate | Acetic acid, ethyl ether; ethyl ether of acetic acid | 0.2 | s.-t. | 2 |
| 3.3.2.1.1.1.1.2. Containing double bonds | | | | |

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| Cis-8-dodecynyl-acetate | Acetic acid, Z-dodec-8-enylic ether; Z-dodec-8-enylic ether of acetic acid; Denacil | 0.00001 | org. odr. | 4 |
| Vinyl acetate | Acetic acid, vinyl ether; vinyl ether of acetic acid | 0.2 | s.-t. | 2 |
| 3.3.2.1.1.1.1.3. of polyatomic alcohols | | | | |
| 3.3.2.1.1.1.1.4. alcohols containing hydroxy, oxy and oxo groups | | 0.6 | | |
| Ethylidendiacetate | Acetic acid, 1-acetoxi-ethyl ether; acetoxi-ethyl ether of acetic acid | 0.6 | s.-t. | 2 |
| 3.3.2.1.1.1.2. of halogen substituted compounds | | | | |
| 2,4,5-Trichlorphenoxiethyl-alpha,alpha-dichlorpropionate | 2,2-dichlor-dichlorpropionic acid,2-(2,4,5-trichlorphenoxy)ethyl ether; 2-(2,4,5-trichlorphenoxy)ethyl ether of 2,2-dichlorpropionic acid; | 2.5 | s.-t. | 3 |

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| | pentanate | | | |
| 2,4,5-Trichlorphenoxy ethyltrichloracetate | Acetic acid, trichlor-2-(2,4,5-trichlorphenoxy) ethyl ether; trichlor-2(2,4,5-trichlorphenoxy) ethyl ether of acetic acid; hexanate | 5.0 | s.-t. | 3 |
| 3.3.2.1.1.1.3. Containing hydroxy, oxy and oxo groups | | | | |
| Ethyl ether of lactic acid | 2-hydroxypropanoic acid, ethyl ether | 0.4 | s.-t. | 3 |
| Acetylacetic acid, methyl ether | Methylacetate, methyl ether of acetylacetic acid | 0.5 <1> | s.-t. | 2 |
| Isopropyl ether of lactic acid | 1-hydroxypropanoic acid, 1-methylethyl ether | 1.0 | s.-t. | 3 |
| Acetopropylacetate | Acetic acid, 4-oxopentyl ether 4-oxopentyl ether of acetic acid | 2.8 <1> | s.-t. | 2 |
| 3.3.2.1.1.1.3.1. of halogen substituted compounds | | | | |
| gamma-Chlorocrotyl ether of dichlorphenoxyacetic acid | 4-Chlorbut-2-enyl ether of 2,4-dichlorphenoxyacetic acid; crotylin | 0.02 | org. odr. | 4 |

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| alpha-methylbenzyl ether of 2-chloroacetyl-acetic acid | 2-Chlor-3-oxooleic acid, 1-phenyl-ethyl ether | 0.15 | s.-t. | 2 |
| Octyl ether of 2,4-dichlorophenoxy-acetic acid | 2,4-dichlorophenoxy-acetic acid, octyl ether | 0.2 | org. odr. | 3 |
| Butyl ether of 2,4-dichlorophenoxy-acetic acid | 2,4-dichlorophenoxy-acetic acid, butyl ether; butyl ether 2,4-D; 2,4-DB | 0.5 | org. odr. | 3 |
| 3.3.2.1.1.2. Containing double or triple bonds | | | | |
| 3.3.2.1.1.2.1. of monohydroxy alcohols | | | | |
| Ethyl acrylate | Acrylic acid, ethyl ether; ethyl ether of acrylic acid | 0.005 | org. odr. | 4 |
| Ethyl ether of 3,3-dimethyl-4,6,6-trichlor-5-hexenic acid | 3,3-dimethyl-4,6,6-trichlor-5-hexenic acid ethyl ether | 0.008 | org. odr. | 3 |
| Butyl acrylate | Acrylic acid, butyl ether; butyl ether of acrylic acid | 0.01 | org. flavr. | 4 |
| Methylmetacrylate | 2-methyl-2-propenic acid, methyl ether; | 0.01 | s.-t. | 2 |

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| | methyl ether of metacrylic acid | | | |
| Butyl ether of metacrylic acid | Metacrylic acid, butyl ether | 0.02 | org. odr. | 4 |
| Methyl acrylate | Acrylic acid, methyl ether; methyl ether of acrylic acid | 0.02 | org. odr. | 4 |
| Ethyl ether of beta, beta-dimethyl acrylic acid | Ethyl ether of 3-methyl-but-2-enic acid | 0.4 | org. odr. | 3 |
| 3.3.2.1.1.2.2. of polyatomic alcohols | | | | |
| Monometacrylic ether of ethylene glycol | Metacrylic acid, 2-hydroxyethyl ether | 0.03 | s.-t. | 4 |
| 3.3.2.1.2. Closed-chain | | | | |
| 3.3.2.1.2.1. Alicyclic | | | | |
| Methyl ether of 2,2-dimethyl-3-propenyl-1-cyclopropanecarbonic acid | 2,2-dimethyl-3-(2-methyl-prop-1-enyl)-cyclopropan-1-carbonic, methyl ether; methyl ether of chrisanthemic acid; methylchrisantemate | 0.61 | org. odr. | 4 |
| 3.3.2.1.2.1.1. Containing oxogroups | | | | |

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| 3.3.2.1.2.2. Aromatic | | | | |
| Methyl benzoate | Benzoic acid, methyl ether; methyl ether of benzoic acid, neobon oil | 0.05 | org. flavr. | 4 |
| n-toluic acid, methyl ether | 4-methyl-benzoic acid methyl ether; methyl ether of n-toluic acid | 0.05 | org. flavr. | 4 |
| 3.3.2.1.2.2.1. Containing aromatic substitute in alcohol | | | | |
| 3.3.2.2. Compound ethers of dibasic acids | | | | |
| 3.3.2.2.1. Aliphatic | | | | |
| 3.3.2.2.1.1. Maximum | | | | |
| 3.3.2.2.1.1.1. Maximum aliphatic alcohols | | | | |
| 3.3.2.2.1.1.2. Olefinic alcohols | | | | |
| 3.3.2.2.1.2. contain- ing double or triple bonds | | 1.0 | | |
| Diethyl ether of maleinic acid | Maleinic acid, diethyl ether | 1.0 | s.-t. | 2 |
| 3.3.2.2.2. Aromatic | | | | |
| Dimethylphthalate | Phthalic acid, dimethyl ether; dimethyl ether of phthalic acid | 0.3 | s.-t. | 3 |
| Diethyl ether of tetrachlortere- | Tetrachlor-terephthalic | 1.0 | s.-t. | 3 |

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| phthalic acid | acid, dimethyl ether; dacthal W-75; chlorthalme- thyl | | | |
| Dimethylterephthalate | Terephthalic acid, dimethyl ether; dimethyl ether of terephthalic acid | 1.5 | org. odr. | 4 |
| 3.3.3. Anhydrides and acid halides | | | | |
| Diacyl chloride of terephthalic acid | Terephthalic acid, dichlor- anhydride; terephtha- loylchloride ; 1,4-ben- zenedicarbo- nyldichlo- ride | 0.02 | org. odr. | 4 |
| Diacyl chloride of 2,3,5,6- tetrachlorterephtha- acid | 2,3,5,6- tetrachlor- terephthalic acid, di- chloranhyd- ride; 2,3,5, 6-tetrachlor terephtha- loyl di- chloride; 2,3,5,6-tet- rachlor-1,4- benzendicar- bonyl- dichloride | 0.02 | org. odr. | 4 |
| Diacyl chloride of isophthalic acid | Isophthalic acid, dichloranhyd ride; isoph- thaloylchlo- ride; 1,3- benzenedicar bonyl- dichloride | 0.08 | org. odr. | 4 |

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| 4. Nitrogen containing compounds | | | | |
| 4.1. Amines and their salts | | | | |
| 4.1.1. Primary | | | | |
| 4.1.1.1. Containing one aminogroup | | | | |
| 4.1.1.1.1. Aliphatic | | | | |
| 4.1.1.1.1.1. Containing only maximum bonds | | | | |
| Amines C16 - C20 | | 0.03 | org. odr. | 4 |
| Amines C10 - C15 | | 0.04 | org. odr. | 4 |
| Monoisopropanolamine | 2-Methyl-1-propanamine | 0.04 | org. flavr. | 3 |
| Amines C7 - C9 | | 0.1 | org. odr. | 3 |
| Monopropylamine | Propylamine | 0.5 | org. odr. | 3 |
| Monoethylamine | Ethylamine | 0.5 | org. odr. | 3 |
| t-Butylamine | | 1.0 | s.-t. | 3 |
| Monomethylamine | Methylamine | 1.0 | s.-t. | 3 |
| Isopropylamine | | 2.0 | s.-t. | 3 |
| Monobutylamine | Butylamine | 4.0 | org. odr. | 3 |
| 4.1.1.1.1.1.1. Containing oxy-, oxo- and carboxygroups | | | | |
| Isopropanolamine | 1-Amino-2-hydroxy-propane | 0.3 | s.-t. | 2 |
| Monoethanolamine | 2-Amino-ethanol | 0.5 | s.-t. | 2 |
| 4.1.1.1.1.1.2. Containing olefinic bonds | | | | |
| Monoallylamine | Allylamine | 0.005 | s.-t. | 2 |
| 4.1.1.1.1.1.2.1. Containing oxy-, oxo-, hydroxy and carboxygroups | | | | |
| Vinyl ether of monoethanolamine | 2-(Ethenyl-oxy)ethanamine, 1-vinyl oxy-2-amino-ethane | 0.006 | org. odr. | 3 |
| 4.1.1.1.1.1.2.2. Acid amides | | | | |

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| Acrylamide | Propenamide, Acrylic acid, amide | 0.01 | s.-t. | 2 |
| Metacrylamide | Metacrylic acid, amide | 0.1 | s.-t. | 2 |
| Methylolmetacrylamide | 4-Hydroxy-2-methylbuten-2-nic acid, amide | 0.1 | s.-t. | 2 |
| N,N-Dimethylaminomethylacrylamide | KF-6 | 2.0 | s.-t. | 2 |
| 4.1.1.1.2. Closed-chain | | | | |
| 4.1.1.1.2.1. Alicyclic | | | | |
| 4.1.1.1.2.2. Aromatic | | | | |
| 4.1.1.1.2.2.1. Mononucleate | | | | |
| Alkylaniline | | 0.003 | s.-t. | 2 |
| 2,4,6-Trimethylaniline | 2,4,6-Trime-thylaniline, mesidine | 0.01 | s.-t. | 2 |
| Aniline | Phenylamine, aminobenzene | 0.1 | s.-t. | 2 |
| n-Butylaniline | n-Aminobu-thylbenzene | 0.4 | org. odr. | 3 |
| m-Toluidine | 3-Methyl-aniline | 0.6 | s.-t. | 2 |
| n-Toluidine | 4-Methyl-aniline, m-aminomethyl-benzene | 0.6 | org. odr. | 3 |
| 4.1.1.1.2.2.1.1. Halogen substituted | | | | |
| Dichloraniline | Dichlorben-benzenamine | 0.05 | org. | 3 |
| Bromotoluene | Bromotolue-dine(mix of o,m,n-isomers) | 0.05 <1> | org. odr. | 4 |

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| m-Trifluoromethyl-aniline | 3-(Trifluoromethyl)benzenamine, 3-aminobenzotri-fluoride | 0.02 | s.-t. | 2 |
| m-Chloraniline | 3-Chlorobenzenamine | 0.2 | s.-t. | 2 |
| n-Chloraniline | 4-Chlorobenzenamine | 0.2 | s.-t. | 2 |
| 2,4,6-Tri-chloroaniline | 2,4,6-Tri-chlorobenzenamine | 0.8 | org. flavr. | 3 |
| 2,4,5-Tri-chloroaniline | 2,4,5-Tri-chlorobenzenamine | 1.0 | org. film | 4 |
| 4.1.1.1.2.2.1.2. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| o-Aminophenol | 1-Amino-2-hydroxybenzene, o-hydroxy-aniline | 0.01 | org. clr. | 4 |
| n-Anisidine | 4-Methoxy-aniline | 0.02 | s.-t. | 2 |
| o-Anisidine | 2-Methoxy-aniline | 0.02 | s.-t. | 2 |
| n-Phenetidine | 4-Ethoxy-aniline, amino-phenetol | 0.02 | s.-t. | 2 |
| n-Aminophenol | | 0.05 | org. clr. | 4 |
| Phenylhydroxylamine | N-Phenylhydroxylamine | 0.1 | s.-t. | 3 |
| m-Aminophenol | 1-Amino-3-hydroxybenzene, hydroxyaniline | 0.1 <1> | org. clr. | 4 |
| 4-Aminobenzoic acid | | 0.1 | s.-t. | 3 |
| 5-Aminosalicylic acid | 5-Amino-2-hydroxy- | 0.5 | org. clr. | 4 |

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| | benzoic acid | | | |
| 3-Aminobenzoic acid | | 10.0 | org. clr. | 4 |
| 4.1.1.1.2.2.1.2.1. Halogen substituted | | | | |
| 4-Amino-3-chlorphenol | | 0.1 | org. clr. | 4 |
| 4.1.1.1.2.2.1.3. Acid amides | | | | |
| Benzamide | | 0.2 <1> | s.-t. | 3 |
| 4.1.1.1.2.2.2. Aromatic condensated | | | | |
| 1-Aminoanthraquinone | | 10.0 | s.-t. | 2 |
| 4.1.1.2. Containig two or more aminogroups | | | | |
| 4.1.1.2.1. Aliphatic | | | | |
| 4.1.1.2.1.1. Containing only maximum bonds | | | | |
| Hexamethylenediamine | 1,6-Diamino-hexane | 0.01 | s.-t. | 2 |
| Hydrazine | | 0.01 | s.-t. | 2 |
| 1,12-Dodecamethylendiamine | 1,12-Dodecan diamine, 1,12-diamino dodecane | 0.05 | s.-t. | 3 |
| Ethylendiamine | 1,2-Diamino-ethane | 0.02 | org. odr. | 4 |
| 4.1.1.2.1.1.1. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| Tetraoxypropylethy-lendiamine | Lapromol 294 | 2.0 | s.-t. | 2 |
| 4.1.1.2.1.1.2. Acid amides | | | | |
| 4.1.1.2.1.2. Containing olefinic bonds | | | | |
| Diallylamine | | 0.01 | s.-t. | 2 |
| Alkylpropylendiamine | | 0.16 | org. odr. | 4 |
| 4.1.1.2.2. Aromatic | | | | |
| 4.1.1.2.2.1. Mononucleate | | | | |
| o-Phenylendiamine | 1,2-Diamino- | 0.01 | org. clr. | 3 |

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| | benzene, phenylen-1,2-diamine | | | |
| Phenylhydrazine | | 0.01 | s.-t. | 3 |
| 4,4'- Diamino-diphenyl ether | 4,4'-Oxy-bisbenzen-amine | 0.03 | s.-t. | 2 |
| m,n-Phenylendiamine | Diaminobenzene, phenylendiamine | 0.1 | s.-t. | 2 |
| 4.1.1.2.2.2. Condensated multinucleate | | | | |
| 1,4-Diamino-anthraquinone | 1,4-Diamino-9,10-anthracendione | 0.02 | org. clr. | 3 |
| 1.5-Diamino-anthraquinone | 1.5-Diamino-9,10-anthracendione | 0.2 | org. clr. | 4 |
| 4.1.2. Secondary | | | | |
| 4.1.2.1. Containing only aliphatic substitutes | | | | |
| Diisobutylamine | Bis(2-methylpropyl)-amine, 2-methyl-M-(2-methylpropyl)-1-propan-amine | 0.07 | org. flavr. | 4 |
| Dimethylamine | | 0.1 | s.-t. | 2 |
| Isopropyloctadecyl-amine | N-Isopropyl-octadecyl-amine | 0.1 | org. film | 4 |
| Diethyltri-amine | N-(2-aminoethyl)-1,2-ethanedi-amine, 2,2'-diaminodiethyl-amine | 0.2 | org. odr. | 4 |
| Dipropylamine | N-propyl-1-propanamine | 0.5 | org. flavr. | 3 |
| Diisopropylamine | M-Isopropyl-1-isopropan-amine | 0.5 | s.-t. | 3 |

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| Ethylbutylamine | N-Ethyl-1-butanamine | 0.5 | org. flavr. | 3 |
| Dibutylamine | N-Butyl-1-butanamine | 1.0 | org. odr. | 3 |
| Diethylamine | | 2.0 | s.-t. | 3 |
| 4.1.2.1.1. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| Diethanolamine | | 0.8 | org. flavr. | 4 |
| 4.1.2.1.2. Oxymes | | | | |
| Acetoxyme | | 8.0 | s.-t. | 2 |
| 4.1.2.1.3. Hydroxamic acids | | | | |
| 4.1.2.2. Containing closed chain substitutes | | | | |
| 4.1.2.2.1. Containing alicyclic substitutes | | | | |
| N-Ethylcyclohexylamine | | 0.1 | s.-t. | 4 |
| 4.1.2.2.1.1. Urea derivatives with one alicyclic substitute | | | | |
| 4.1.2.2.2. Containing mononucleate aromatic substitutes | | | | |
| 4-Aminodiphenylamine | N-Phenyl-1,4 benzenedi-amine, N-phenyl-n-phenylen-diamine | 0.005 | s.-t. | 2 |
| Diphenylamine | N-Phenylbenzenamine | 0.05 | org. odr. | 3 |
| N-Methylaniline | | 0.3 | org. odr. | 2 |
| N-Ethyl-o-toluidine | N-Ethyl-2-methyl-aniline | 0.3 | org. odr. | 3 |
| N-Ethylmetatoluidine | 3-Methyl-N-ethylaniline | 0.6 | s.-t. | 2 |
| N-Ethylaniline | N-Ethylbenzenamine | 1.5 | org. odr. | 3 |
| 4.1.2.2.2.1. Containing hydroxy, oxy, oxo and carboxygroups | | | | |

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| 4-Amino-2-(2-hydroxyethyl)-N-ethylaniline sulphite | | 0.2 | org. odr. | 3 |
| n-Acetaminophenol | Acetic acid, (4-hydroxyphenyl)-amide; paracetamol; 4-acetamidophenol | 1.0 | org. flavr. | 3 |
| N-Acetyl-2-aminophenol | | 2.5 | org. clr. | 4 |
| 4.1.2.2.2.2. Oxymes | | | | |
| Cyanbenzaldehyde oxyme, sodium salt | | 0.03 | org. odr. | 4 |
| n-Quinone-dioxyme | 2,5-Cyclohexanediene-1,4-dione dioxyme | 0.1 | s.-t. | 3 |
| Cyclohexane-oxyme | | 1.0 | s.-t. | 2 |
| 4.1.2.2.2.3. Acid amides | | | | |
| 3-Chlor-2,4-dimethyl-valerateaniline | 2-Methyl 2-methylpentanic acid, 4-methyl-3-chloranilide; solan | 0.1 | org. odr. | 4 |
| Anilide of salicylic acid | | 2.5 | org. odr. | 3 |
| 4.1.2.2.2.4. Urea derivatives with one aromatic substitute | | | | |
| m-Trifluoromethylphenylurea | 1-(3-Trifluoromethylphenyl) urea | 0.03 | org. flavr. | 4 |
| 4-Chlor-2-butynyl-N-(3-chlorphenyl)-carbamate | 4-Chlorphenylcarbamic acid, 4-chloro-but-2-ynic ether, | 0.03 | org. odr. | 4 |

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| | carbine | | | |
| 3-Methylphenyl-N-methylcarbamate | Methylcarbamic acid, methylphenyl ether; dicresyl | 0.1 | org. odr. | 3 |
| Isopropylphenyl-carbamate | Phenyl-carbamic acid, isopropyl ether | 0.2 | org. odr. | 4 |
| Isopropylchlorophenyl carbamate | 3-chlorophenyl carbamic acid, isopropyl ether | 1.0 | org. odr. | 4 |
| Oxyphenylmethylurea | 1-Hydroxy-3-methyl-1-phenyl urea; meturine | 1.0 | s.-t. | 3 |
| 3-Methoxycarbamido-phenyl-N-phenylcarbamate | 3-tolylcarbamic acid, 3-(N-methoxycarbonylamino)phenyl ether; Phenmedipham | 2.0 | s.-t. | 3 |
| 4.1.2.2.3. Containing polynuclear aromatic substitutes | | | | |
| 1-Chlor-4-benzoylaminoanthraquinone | | 2.5 | s.-t. | 3 |
| 4.1.2.2.3.1. Urea derivates with condensated aromatic substitute | | | | |
| 1-Naphthyl-N-methylcarbamate | Methylcarbamic acid, naphth-1-yl ether; sevin | 0.1 | org. odr. | 4 |
| 4.1.3. Tertiary | | | | |
| 4.1.3.1. Containing only aliphatic substitutes | | | | |

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| Triallylamine | | 0.01 | s.-t. | 2 |
| 1-Butylbiguanidine hydrochloride | Glybutide | 0.01 <1> | s.-t. | 2 |
| Triisooctylamine | N,N-Diisooctyl isooctylamine | 0.025 | s.-t. | 2 |
| Trimethylamine | | 0.05 | org. odr. | 4 |
| Trialkylamine C7 - C9 | | 0.1 | s.-t. | 3 |
| Alkyldimethylamine | | 0.2 | s.-t. | 3 |
| N,N'-Diethylguanidine, muriatic | 1,2-Diethylguanidine monohydrochloride | 0.8 | s.-t. | 3 |
| Tributylamine | | 0.9 | org. odr. | 3 |
| Triethylamine | | 2.0 | s.-t. | 2 |
| 4.1.3.1.1. Nitriles | | | | |
| Malononitrile | Propanedinitrile, dicyano methane | 0.02 | s.-t. | 2 |
| Acetonecyanohydrin | 2-hydroxy-2-methylpropanoic acid, nitrile; 2-hydroxy-2-methylpropanonitrile, nitrile of hydroxyisobutyric acid | 0.035 | s.-t. | 2 |
| Alkylaminopropionitrile C17 - C20 | | 0,05 | org. foam | 4 |
| Dinitrile of adipic acid | | 0.1 | s.-t. | 2 |
| Allyl cyanide | But-3-enoic acid, nitrile | 0.1 | s.-t. | 2 |
| Isocrotononitrile | 2-Methyl-2-propen- | 0.1 | s.-t. | 2 |

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| | nitrile | | | |
| Crotonitrile | But-2-enic acid, nitrile | 0.1 | s.-t. | 2 |
| Succinonitrile | Butanedinitrile | 0.2 | s.-t. | 2 |
| Acetonitrile | Acetic acid, nitrile | 0.7 | org. odr. | 3 |
| Calcium cyanamide | Carbamic acid nitrile, compound with calcium | 1.0 | s.-t. | 3 |
| Acrylic acid nitrile | | 2.0 | s.-t. | 2 |
| Dicyandiamide | Cyanoguanidine | 10.0 | org. flavr. | 4 |
| 4.1.3.1.2. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| Triisopropanolamine | Tripropylamine | 0.5 | s.-t. | 2 |
| Triethanolamine | | 1.0 | org. flavr. | 4 |
| Ethyl ether of N-benzoil-N-(3,4-dichlorophenyl)-2-aminopropionic acid | Ethyl-N-benzoil-N-(3,4-dichlorophenyl)alaninate, suffix | 1.0 | s.-t. | 2 |
| Methyldiethanolamine | Bis(2-hydroxyethyl)methylamine, 2,2-(M-methylamino) diethanol | 1.0 | s.-t. | 2 |
| 4.1.3.1.3. Amides | | | | |
| Dimethylacetoamide | | 0.4 | s.-t. | 2 |
| Diethylamide 2-(alpha naphthoxy) propionic acid | N,N-Diethyl-2-(1-naphthalenyloxy)propanamide | 1.0 | s.-t. | 2 |

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| 4.1.3.1.4. Urea derivatives with several aliphatic substitutes | | | | |
| N,N-Dimethylurea | 1,3-Dimethyl urea | 1.0 | s.-t. | 2 |
| N,N-Diethyl carbamylchloride | | 6.0 | s.-t. | 2 |
| 4.1.3.2. Containing closed chain substitutes | | | | |
| 4.1.3.2.1. Urea derivatives with alicyclic substitutes | | | | |
| 3-(Hexahydro-4,7-methanindan-5-yl)-1,1-dimethylurea | Herban | 2.0 | s.-t. | 2 |
| 4.1.3.2.2. Containing aromatic substitutes | | | | |
| N,N-Dimethyl-n phenylendiamin-sulphate | CPV, 1,4-amonodiethylanilinsulphate | 0.1 | s.-t. | 2 |
| N,N-Diethylaniline | N,N-Diethylbenzenamine | 0.15 | org. clr. | 3 |
| Alkylbenzyldimethylammonium chloride C10-C16 | | 0.3 | org. foam | 3 |
| Alkylbenzyldimethylammonium chloride C17-C20 | | 0.5 | org. foam | 3 |
| N-(C7 - C9)Alkyl-N-phenyl-n-phenylendiamine | C-789 product | 0.9 <1> | org. clr. | 3 |
| Ethylbenzylaniline | N-Phenyl-N-ethylbenzenmethanamine | 4.0 | s.-t. | 2 |
| 4.1.3.2.2.1. Nitriles, isonitriles | | | | |
| Benzyl cyanide | Isocyanomethylbenzene | 0.03 | org. odr. | 4 |
| Dinitrile of isophthalic acid | 1,3-Benzene-dicarbonitrile, isophthalonitrile, 1,3-dicyanobenzene | 5.0 | s.-t. | 3 |

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| 4.1.3.2.2.2. Amides | | | | |
| 4.1.3.2.2.3. Urea derivatives with one or several aromatic substitutes | | | | |
| Diphenylurea | N,N-Diphenyl urea, carbanilide | 0.2 | org. odr. | 4 |
| N-Trifluoromethyl-phenyl-N',N'-dimethylurea | 1,1-Dimethyl 3-(3-trifluoromethyl-phenyl)urea, cotoran | 0.3 | org. film | 4 |
| Diethylphenylurea | Centralite | 0.5 | org. flavr. | 4 |
| N'-(3,4-Dichlorophenyl)-N,N-dimethylurea | 1,1-Dimethyl 3-(3,4-dichlorophenyl) urea, diuron | 1.0 | org. odr. | 4 |
| 4.1.4. Quaternary ammonium salts | | | | |
| Nitrate of methyltri-alkyl ammonium | | 0.01 | s.-t. | 2 |
| Alkyltrimethyl-ammonium chloride | | 0.2 | s.-t. | 2 |
| Chlorcholine chloride | N,N,N-Trime-thyl-N-(2-chloroethyl) ammonium chloride | 0.2 | s.-t. | 2 |
| 4.2. Oxygen and nitrogen containing compounds | | | | |
| 4.2.1. Nitro- and nitroso compounds | | | | |
| 4.2.1.1. Aliphatic | | | | |
| Nitromethane | | 0.005 | org. odr. | 4 |
| Trinitromethane | Nitroform | 0.01 | org. clr. | 3 |
| Tetranitromethane | | 0.5 | org. odr. | 4 |
| Nitropropane | | 1.0 | s.-t. | 3 |
| Nitroethane | | 1.0 | s.-t. | 2 |
| 4.2.1.1.1. Containing hydroxy, oxy, oxo and carboxygroups | | | | |

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| Dinitrodiethyleneglycol | Dihydroxyethyl ether dinitrate, diethyleneglycol dinitrate | 1.0 | s.-t. | 3 |
| Dinitrotriethyleneglycol | | 1.0 | s.-t. | 3 |
| 4.2.2. Closed-chain | | | | |
| 4.2.2.1. Alicyclic | | | | |
| Chloronitrosocyclohexane | 1-Nitroso-1-chlorocyclohexane | 0.005 | org. odr. | 4 |
| Nitrocyclohexane | | 0.1 | s.-t. | 2 |
| 4.2.1.2.2. Aromatic | | | | |
| 4.2.1.2.2.1. Mononucleate | | | | |
| Nitrobenzene | | 0.2 | s.-t. | 3 |
| Trinitrobenzene | | 0.4 | s.-t. | 2 |
| Dinitrobenzene | | 0.5 | org. odr. | 4 |
| 2,4-Dinitrotoluene | | 0.5 | s.-t. | 2 |
| 4.2.1.2.2.1.1. Halogen substituted | | | | |
| n-Trifluoromethylnitrobenzene | 1-Nitro-3-trifluoromethylbenzene | 0.01 | org. odr. | 3 |
| Nitrochlorobenzene | Nitrochlorobenzene (mix of 2,3,4 isomers) | 0.05 | s.-t. | 3 |
| Nitroisophenol | | 0.1 | org. clr. | 3 |
| 2,5-Dichloronitrobenzene | 1,4-Dichloro-2-nitrobenzene | 0.1 | s.-t. | 2 |
| 3,4-Dichloronitrobenzene | 4-Nitro-1,2-dichlorobenzene | 0.1 | s.-t. | 3 |
| Dinitrochlorobenzene | 2,4-Dinitro- | 0.5 | org. odr. | 3 |

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| | 1-benzene chloride | | | |
| 4.2.1.2.2.1.2. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| n-Nitrophenetol | 4-Nitro-ethoxy-benzene | 0.002 | s.-t. | 2 |
| n-Nitrophenetol | 4-Nitro-phenol | 0.02 0.02 | s.-t. s.-t. | 2 2 |
| 2-sec-Butyl-4,6-dinitrophenyl-3,3-dimethylacrylate | 2-(1-Methyl-propyl)-4,6-dinitrophenyl-3-methyl-2-butenolate, morocide, acricide, endosane, 2-sec-butyl-4,6-dinitrophenyl-3-methyl-crotonate | 0.03 | s.-t. | 2 |
| 2,4-Dinitrophenol | | 0.03 | s.-t. | 3 |
| 2-Methyl-4,6-dinitrophenol | | 0.05 | s.-t. | 2 |
| m-Nitrophenetol | 3-Nitro-phenol | 0.06 | s.-t. | 2 |
| o-Nitrophenetol | 2-Nitro-phenol | 0.06 | s.-t. | 2 |
| n-Nitroanisole | 4-Nitromethoxybenzene | 0.1 | org. flavr. | 3 |
| 2-(1-Methyl-propyl)-4,6-dinitrophenol | Dinoseb | 0.1 | org. clr. | 4 |
| m-Nitrobenzoic acid | 3-Nitro-benzoic acid | 0.1 | org. clr. | 4 |
| n- Nitrobenzoic acid | 4-Nitro-benzoic acid | 0.1 | s.-t. | 3 |
| Methylethyl-[2-(1- | 2-sec-butyl- | 0.2 | org. | 4 |

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| ethylmethylpropyl)- 4,6-dinitrophenyl] carbonate | 4,6-dinitro- phenylic acid, isopropyl ether; dinobutone; citazole; acrex | | film | |
| o-Nitroanisole | 2-Nitroani- sole | 0.3 | org. flavr. | 3 |
| 2,4,6- Trinitrophenol | Picronitric acid | 0.5 | org. clr. | 3 |
| 2-[(n-Nitrope- nyl)acetylami- no]ethan-1-ol | Oxyacetyl- amine | 1.0 | org. odr. | 4 |
| 4.2.1.2.2.1.2.1. Halogen substituted | | | | |
| n-Nitrophenylchloro- metincarbinol | 4-Nitro- alpha- chloromethyl benzenmetha- nol; [1-(4- nitro- phenyl)] -2-chloretha- -1-ol | 0.2 | org. odr. | 4 |
| 3-Nitro-4-chloro- benzoic acid | | 0.25 | org. flavr. | 3 |
| 5-Nitro-2-chloro- benzoic acid | | 0.3 | org. flavr. | 4 |
| 2,5-Dichlor-3- nitrobenzoic acid | | 2.0 | s.-t. | 2 |
| 2,4-Dichlorophenyl- 4-nitrophenylic ether | 2,4-Dichlor- 1-(4-nitro- phenoxy) benzene, nitro- chlor, Tokkorn | 4.0 | s.-t. | 2 |
| 4.2.1.2.2.1.3. Containing amino-, imino- and diazogroups | | | | |
| 4-Nitro-N,N- diethylaniline | | 0.002 | org. clr. | 3 |
| 2-Nitroaniline | o-Nitroani- | 0.01 | org. clr. | 3 |

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| | line | | | |
| N-Nitroisodiphenylamine | Diphenyl-nitrosamine | 0.01 | s.-t. | 2 |
| 2,4-Dinitro-2,4-diazopentane | N,N'-Dimethyl-N,N-dinitromethandiamine | 0.02 | s.-t. | 2 |
| 4-Nitroaniline | n-Nitroaniline, 4-nitrobenzenamine | 0.05 | s.-t. | 3 |
| Dinitroaniline | Dinitrobenzenamine | 0.05 | org. clr. | 4 |
| 3-Nitroaniline | 3-Nitrobenzenamine, m-nitroaniline | 0.15 | org. clr. | 3 |
| Indotoluidine | N-(4-Amino-3-methylphenyl)-n-benzoquinonimine | 1.0 | s.-t. | 2 |
| 4.2.1.2.2.1.3.1. Halogen substituted | | | | |
| 4-Chloro-2-nitroaniline | 4-Chloro-2-nitrobenzenamine | 0.025 | org. clr. | 3 |
| 2,6-Dichloro-4-nitroaniline | 2,6-Dichloro-4-nitrobenzenamine, dichlorane, botran | 0.1 | org. | 3 |
| 3,5-Dinitro-4-diethylaminobenzotrifluoride | Nitrofor | 1.0 | org. odr. | 4 |
| 3,5-Dinitro-4-dipropylaminobenzotrifluoride | 2,6-Dinitro-N,N-dipropyl-4-trifluoromethyl-aniline, Treflan | 1.0 | org. odr. | 4 |
| 4.2.1.2.2.1.3.2. Containing hydroxy, oxy, oxo and carboxygroups | | | | |
| 2,4,4-Trinitrobenzanilide | 2,4,6-Trinitro-6-trinitro- | 0.02 | s.-t. | 2 |

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| | benzoic acid, anilide | | | |
| n-Nitrophenylamino-ethanol | 2-[(4-nitrophenyl)amino]ethanol, oxy-amine | 0.5 | org. odr. | 4 |
| 4.2.1.2.2.2. Aromatic condensated | | | | |
| Dinitronaphthalene | | 1.0 | org. clr. | 4 |
| 1-Nitroanthraquinon-2-carboxylic acid | 9,10-dihydro-1-nitro-9,10-dioxo-2-anthracene acid | 2.5 | s.-t. | 3 |
| 4.2.2. Ethers and salts of nitric and nitrous acids | | | | |
| Butylnitrite | Nitrous acid, butyl ether | 0.05 | org. odr. | 4 |
| 1-Nitroguanidine | | 0.1 | s.-t. | 2 |
| 5. Sulphur-containing compounds | | | | |
| 5.1. Thiocompounds | | | | |
| 5.1.1. Containing C-S-H group | | | | |
| Methylmercaptan | | 0.0002 | org. odr. | 4 |
| Allylmercaptan | | 0.0002 | org. odr. | 3 |
| beta-Mercaptodiethylamine | 2-(N,N-Diethylamino)-ethanethiol | 0.1 | org. odr. | 4 |
| 5.1.2. Containing C-S-C group | | | | |
| Dimethyl sulphide | | 0.01 | org. odr. | 4 |
| 3-Methyl-4-methylthiophenol | Methylthiomethylphenol, 3-methyl-4-thioanisole | 0.01 | org. flavr. | 4 |
| 2-Methylthio-O-methylcarbomol-butanoxy-3 | 3-Methylthio-2-butanol-O-(methylami- | 0.1 | org. odr. | 3 |

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| | nocarbonyl) oxyme, Dravine 755 | | | |
| 4-Chlorophenyl-2,4,5-trichlorophenyl-sulphide | 1,2,4-Tri-chloro-5-[4-(chloro-phenyl)thio]benzen-tetrazole, animert | 0.2 | org. film | 4 |
| Divinylsulphide | Vinylsul-phide, 1,1-thio-bisethene | 0.5 | org. odr. | 3 |
| 5.1.3. Containing C-S-S-C group | | | | |
| Dimethyldisulphide | | 0.04 | org. odr. | 3 |
| 5.1.4. Containing C=S group | | | | |
| Carbon disulphide | | 1.0 | org. odr. | 4 |
| 5.1.4.1. Thiourea derivatives | | | | |
| S-Propyl-N-ethyl-N-butylthiocarbamate | Butyl (ethyl)thio carbamic acid, S-propylyc ether; tillam | 0.01 | org. odr. | 3 |
| Thiourea | Thiocarbami-de, diamide of thiocarba mic acid | 0.03 | s.-t. | 2 |
| S-(2,3-Dichlorallyl)-N,N-diisopropylthio-carbamate | Diisopropyl-thiocarbamic acid, S-(2,3-di-chloroprop-2-enylic) ether; avadex | 0.03 | org. odr. | 4 |
| S-Ethyl-N,N'-dipropylthiocarbamate | Dipropylthio carbamic acid, S-ethyl ether; eptam | 0.1 | org. odr. | 3 |

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| Amidine-thio acetic acid | Carboxymethylisothiourea | 0.4 | s.-t. | 2 |
| 1,2-Bis-methoxycarbonylthioureidobenzen | 1,2-phenylen-bis (imino-(carbonothioyl)biscarbamic acid, diethyl ether; topsin; nemafax; thiophanate | 0.5 | org. flavr. | 3 |
| 5.1.4.2. Dithiocarbarnic acid derivatives | | | | |
| Tetraethyl thiuram disulphide | N,N,N',N'-Tetraethyl thiuram disulphide, thiuram E | None | org. odr. | 3 |
| N- methyl-dithiocarbamic acid, N-methylamine salt | | 0.02 | org. odr. | 3 |
| Sodium methyldithiocarbamate | Methyldithiocarbamic acid, sodium salt; carbathion | 0.02 | org. odr. | 3 |
| Ammonium ethylenbisthio carbamate | 1,2-ethylenbisthio-carbamic acid, diammonium salt | 0.04 | org. odr. | 3 |
| S-Ethyl-N-ethyl-N-cyclohexylthiocarbamate | Ronite, cycloate | 0.2 | s.-t. | 3 |
| Zink ethylenbisdithiocarbamate | N,N'-ethylenbisdithiocarbamic acid, zink salt; zineb | 0.3 | org. МУТН. | 3 |
| Ammonium dimethyldithiocarbamate | Dimethyldithiocarbamic | 0.5 | s.-t. | 3 |

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| | acid, ammonium salt | | | |
| Tetramethylthiuram disulphide | Tetramethylthiuram disulphide, thiuram D | 1.0 | s.-t. | 2 |
| 5.1.4.3. Xanthogenates | | | | |
| Butylxanthogenate | Thiolthiocarbonic acid, butyl ether | 0.001 | org. odr. | 4 |
| Isoamylxanthogenate | Thiolthiocarbonic acid, isoamyl ether; isopentylxanthogenate | 0.005 | org. odr. | 4 |
| Isopropylxanthogenate, salt | Thiolthiocarbonic acid, isopropyl ether; salt | 0.05 | org. odr. | 4 |
| Ethylxanthogenate, salt | Thiolthiocarbonic acid, ethyl ether, salt | 0.1 | org. odr. | 4 |
| 5.1.5. Containing C - N = S group | | | | |
| 5.1.6. Sulphonium salts | | | | |
| (4-Hydroxy-2-methylphenyl) dimethylsulphonium chloride | | 0.007 | org. odr. | 4 |
| 5.2. Compounds containing sulphur directly bonded with oxygen | | | | |
| 5.2.1. Sulphoxydes | | | | |
| 5.2.2. Sulphones | | | | |
| N-n-Butyl-N-(n-methylbenzylsulphonyl)urea | 1-Butyl-1-(n-tolylsulphonyl)-urea, | 0.001 <1> | s.-t. | 1 |

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| | butamide | | | |
| N-Propyl-N'-(n-chlorobenzensulfonyl)urea | 3-Propyl-1-[(n-chloro-nyl)sulphonyl]urea, chloropropamide | 0.001 <1> | s.-t. | 1 |
| 4,4'-Dichlordiphenylsulphone | 1,1'-Sulphonyl-bis(4-chlorobenzene), di-4-chlorophenylsulphone, bis(n-chlorophenyl)sulphone | 0.4 | s.-t. | 2 |
| 4,4'-Diaminodiphenylsulphone | 4.4'-Sulphonyldianiline | 1.0 | s.-t. | 2 |
| 5.2.3. Sulphinic acids and their derivative compounds | | | | |
| n-Toluenesulphonic acid, salt | 4-methylbenzenesulphinic acid, salt | 1.0 | s.-t. | 2 |
| 5.2.4. Sulphonic acids and their derivative compounds | | | | |
| 5.2.4.1. Aliphatic sulphonic acids and their salts | | | | |
| Methyltrialkylammonium methyl sulphate | | 0.01 | s.-t. | 3 |
| Olefin sulphonate C15 - C18 | | 0.2 | s.-t. | 2 |
| olefin sulphonate C12 - C14 | | 0.4 | org. foam | 4 |
| N-methylsulphaminic acid | | 0.4 | s.-t. | 2 |
| Alkylsulphonates | | 0.5 | org. clr. | 4 |
| 5.2.4.2. Aromatic | | | | |
| 5.2.4.2.1. Mononucleate | | | | |
| 5.2.4.2.1.1. Sulphonic acids and their salts containing no substitutes except for alkyl | | | | |

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| Alkylbenzene-sulphonates | Sulphonol chlorine | 0.5 | org. foam | 4 |
| 5.2.4.2.1.1.1. Containing substitutes in the chemical group | | | | |
| 1,4-Bis(4-methyl-2-sulphophenylamino)-5,8-dihydroxyanthraquinone, disodium salt | 2Zh chromic green anthraquinonic coloring agent | 0.01 | org. clr. | 4 |
| 4-nitroanilin-2-sulphonic acid, salt | 4-Nitroaniline-2-sulphonic acid salt | 0.08 | org. clr. | 4 |
| Aminobenzen-3-sulphonic acid | Metalline acid, anilin-m-sulphonic acid | 0.7 | org. clr. | 4 |
| 3-Nitroanilin-4-sulphonic acid | 4-Amino-2-nitrobenzen-sulphonic acid, 3-nitrosulphanilic acid | 0.9 | org. clr. | 4 |
| Sodium n-chlorobenzen-sulphonate | 4-Chlorobenzen-sulphonic acid, sodium salt; Ludigol | 2.0 | s.-t. | 2 |
| 5.2.4.2.1.1.2. Aromatic sulphonic acids ethers | | | | |
| 5.2.4.2.1.1.3. Acid halides of aromatic sulphonic acids | | | | |
| Benzenesulphochloride | Benzenesulphonil-chloride | 0.5 | org. odr. | 4 |
| 5.2.4.2.1.1.4. Amides | | | | |
| n-Butylamide of benzenesulphonic acid | Benzenesulphonic acid, n-butylamide; N-butylbenzen-sulphamide | 0.03 | s.-t. | 2 |

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| Benzensulphamide | Benzensulphonic acid, amide | 6.0 | s.-t. | 3 |
| 5.2.4.2.2. Condensated polynuclear | | | | |
| Bis(n-butyl-anilin) anthraquinone-3,3-disulphonic acid, disodium salt | H ₂ C acidic anthraquinonic green coloring agent | 0.04 | org. clr. | 4 |
| 1,8-diaminonaphthalen-4-sulphonic acid | C-acid | 1.0 | org. odr. | 3 |
| 2-Naphthol-6-sulphonic acid | 6-Hydroxy-2-naphthalen sulphonic acid, beta-naphtholsulphonic acid, Schaffer salt; | 4.0 | s.-t. | 3 |
| 5.3. Ethers and salts of sulphuric and sulphurous acids | | | | |
| 4-Chlorophenyl-4-chlorobenzen-sulphonate | Sulphonate ether | 0.2 | org. flavr. | 4 |
| 2-Aminoethyl ether of sulphuric acid | 2-amino-ethyl sulphuric acid | 0.2 | s.-t. | |
| n-Methylaminophenol sulphate | Metol | 0.3 | org. clr. | 3 |
| Alkylsulphates | | 0.5 | org. foam | 4 |
| Triethanolamine alkylbenzene sulphate | | 1.0 | org. foam | 3 |
| 6. Phosphorus compounds | | | | |
| 6.1. Containing C - P bond | | | | |
| 6.1.1. Phosphines and phosphonium salts | | | | |
| Tris(diethylamino)-2-chlorethyl-phosphine | Dephos | 2.0 | org. odr. | 3 |
| 6.1.2. Tertiary phosphines oxydes | | | | |

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| Triisopentylphosphine oxyde | Tris(3-methylbutyl)-phosphoric acid | 0.3 | s.-t. | 2 |
| Dioctylisopentylphosphine oxyde | (3-Methylbutyl)dioctylphosphine oxyde | 1.0 | s.-t. | 3 |
| 6.1.3. Phosponates | | | | |
| 2-Chlorethylphosphonic acid, bis(2-chlorethyl) ether | Diether of 2-chlorethyl phosphonic acid | 0.2 | s.-t. | 2 |
| Vinylphosphonic acid, bis(beta,beta-хлорэтиловый) ether | O,O-Bis(2-chloroethyl) vinylphosphonate, Vinifos | 0.2 <1> | s.-t. | 2 |
| O,O-Diphenyl-1-hydroxy-2,2,2-trichlorethylphosphonate | | 0.3 | org. foam | 3 |
| O-(2-Chlor-4-methylphenyl) | (4-Methyl-2-chloro-) phenyl | 0.4 | org. odr. | 4 |
| N'-isopropylamido-chloromethylthiophosphonate | N-sec-butylamido-chloromethyl thiophosphonate, Izofos-3 | | | |
| Oxyhexylidenediphosphonate | | 0.5 | s.-t. | 3 |
| Oxyheptylidenediphosphonate | | 0.5 | s.-t. | 3 |
| Oxynonylidenediphosphonate | | 0.5 | s.-t. | 3 |
| Oxyoctylidenediphosphonate | | 0.5 | s.-t. | 3 |
| Oxyethylidendi-phosphonic acid | Hydroxy - ethane-1,1-diphosphonic acid | 0.6 | org. flavr. | 4 |

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| 2-Chlorethylphosphonic acid, 2-chlorethyl ether | Monoether of 2-Chlorethyl phosphonic acid | 1.5 | s.-t. | 3 |
| 2-Chlorethylphosphonic acid | Ethrel, Ethephon, Florel | 4.0 | s.-t. | 2 |
| 2-Hydroxy-1,3-propylendiamine-N,N,N',N'- tetramethylenphosphonic acid sodium salt | DPF-1N | 4.0 | org. flavr. | 4 |
| 6.2. Phosphoric and phosphorous acids derivatives | | | | |
| 6.2.1. Phosphites | | | | |
| Trimethylphosphite | | 0.005 | org. odr. | 4 |
| Triphenylphosphite | O,O,O-Triphenylphosphite | 0.01 | s.-t. | 2 |
| Trimethylphosphite | | 0.02 | org. odr. | 3 |
| 6.2.3. Phosphoric acid amides | | | | |
| 6.2.2. Phosphates | | | | |
| O,O,O-Tricresyl phosphate | Tricresylphosphate | 0.005 | s.-t. | 2 |
| O,O,O-Tributylphosphate | Tributylphosphate | 0.01 | org. flavr. | 4 |
| O,O,O-Trixylenylphosphate | Trixylenylphosphate | 0.05 | org. odr. | 3 |
| O,O-Dimethyl-O-[3-(carb-1-phenylethoxy)propen-2-yl-2-phosphate | 3-dimethoxyphosphoryloxycrotonic acid, 1- phenylethyl ether; Ciodrin | 0.05 | s.-t. | 2 |
| O,O-Dimethyl-O-[1-(2,3,4,5-tetrachlorophenyl)-2-chlorovinyl | Vinylphosphate | 0.2 | org. flavr. | 3 |

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| phosphate | | | | |
| O,O,O-Trimethylphosphate | Trimethylphosphate | 0.3 | org. odr. | 4 |
| 6.2.2.1. Halogen substituted | | | | |
| O,O-Dimethyl-(1-hydroxy-2,2,2-trichlorethyl)phosphonate | Chlorophos | 0.05 | org. odr. | 4 |
| O,O-Dimethyl-O-(2,2-dichlorovinyl)-phosphate | O-(2,2-Dichlorovinyl) O,O-dimethylphosphate, DDVF, dichlorvos | 1.0 | org. odr. | 3 |
| Dichloropropyl (2-ethylhexyl)phosphate | | 6.0 | org. | 4 |
| 6.2.2.2. Thiophosphates | | | | |
| S,S,S-Tributyltrithiophosphate | Butyvos | 0.0003 | org. flavr. | 4 |
| O-Cresyldithiophosphate | Cresylic dithiophosphate | 0.001 | org. odr. | 4 |
| O,O-Dimethyl-S-ethylmercaptoethyl-dithiophosphate | O,O-Dimethyl S-(2-ethylthioethyl) dithiophosphate, M-81 | 0.001 | org. odr. | 4 |
| O,O-Dimethyl-O-(3-methyl-4-methylthiophenyl)-thiophosphate | Thiophosphoric acid, O,O-dimethyl-O-(3-methyl-4-methylthio) phenyl ether; Sulphidovos; Baytex | 0.001 | org. odr. | 4 |
| O-(4-Methylthiophenyl)-O-ethyl-S-propyldithiophosphate | Bolstar, Gelotion, Sulprofos | 0.003 | org. odr. | 4 |
| Bis(2-ethylhexyl)- | Dithiophos- | 0.02 | s.-t. | 2 |

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| dithiophosphoric acid | phoric acid, O,O-bis(2-ethylhexyl) ether | | | |
| O,O-Diethyl-S-carbethoxymethylthiophosphate | Acetophos | 0.03 | org. odr. | 4 |
| O,O-Dimethyl-S-carbethoxymethylthiophosphate | (dimethoxy-(thiophosphorylthio) acetic acid, ethyl ether; Methylacetophos | 0.03 | org. odr. | 4 |
| O,O-Dimethyl-S-(1,2-dicarbethoxyethyl)dithiophosphate | 2-(dimethoxythiophosphorylthio) butandionic acid, diethyl ether; Karbofos | 0.05 | org. odr. | 4 |
| O,O-Diethyl-S-benzylthiophosphate | S-Benzyl-O, O-diethylthio phosphate, Rucid-II | 0,05 | s.-t. | 2 |
| O-Phenyl-ethylthiophosphoric acid, salt | | 0.1 | org. odr. | 4 |
| Dibutyldithiophosphates | Dithiophosphoric acid, O,O-di-butyl ether, salt | 0.1 | s.-t. | 2 |
| Dibutylmonothio-phosphate | | 0.1 | org. odr. | 3 |
| Dimethyldithiophosphoric acid | O,O-dimethyl dithiophosphoric acid | 0.1 | org. odr. | 4 |
| S-(2-Acetamidoethyl)-O,O-dimethyldithiophosphate | Amifos | 0.1 | org. odr. | 4 |

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| Diethyldithio-phosphoric acid | O,O'-Diethyl dithio-phosphoric acid | 0.2 | org. odr. | 4 |
| Diethyldithio-phosphate | Diethyl-dithiophosphoric acid, salt | 0.5 | org. odr. | 3 |
| 6.2.2.2.1. Halogen substituted | | | | |
| O-Methyl-O-ethylchlorthio-phosphate | Diether | 0.002 | org. odr. | 4 |
| O-Phenyl-O-ethylchlorthio-phosphate | | 0.005 | org. odr. | 3 |
| O-(4-Brom-2,5-dichlorophenyl)-O,O-dimethylthiophosphate | Bromophos | 0.01 | org. odr. | 4 |
| Monomethyl-di-chlorthiophosphate | O-Methyl-di-chlorthio-phosphate | 0.01 | s.-t. | 2 |
| Monoethyldichlor-thiophosphate | O-Ethyldi-chlorthio-phosphate | 0.02 | org. odr. | 4 |
| O-(2,4-Dichlorophenyl)-S-propyl-O-ethylthiophosphate | Etaphos, Protiofos, Tokution, Bideron | 0.05 | org. odr. | 3 |
| Diethylchlorthio-phosphate | O,O-Diethyl-chlorthio-phosphate | 0.05 | org. odr. | 4 |
| Dimethylchlorthio-phosphate | O,O-Di-methyl-chlorthio-phosphate | 0.07 | org. odr. | 3 |
| O-Methyl-O-(2,4,5-trichlorophenyl)-O-ethylthiophosphate | Trichlormeta phos-3 | 0.4 | org. odr. | 4 |
| O,O-Dimethyl-O-(2,5-dichlor-4-iodophenyl) thiophosphate | Iodofenfos | 1.0 | org. odr. | 3 |

| 6.2.2.2.2. Nitrogen containing compounds | | | | |
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| O,O-Diethyl-O-(4-nitrophenyl) thiophosphate | O-(4-Nitrophenyl)-O,O-diethylthiophosphate, Thiophos | 0.003 | org. odr. | 4 |
| O,O-Dimethyl-S-(N-methyl-N-formylcarbamoylmethyl)-dithiophosphate | O,O-Dimethyl S-(N-methyl-N-formylaminomethyl)-dithiophosphate, Antio | 0.004 | org. odr. | 4 |
| O,O-Dimethyl-O-(4-nitrophenyl)phosphate | Metaphos | 0.02 | org. odr. | 4 |
| Butylamide of O-ethyl S-phenyldithiophosphoric acid | O-Ethyl-S-phenyl-N-butylamido-dithiophosphate, Fosbutyl | 0.03 | org. odr. | 4 |
| O,O-Dimethyl-S-(N-methylcarbamidomethyl)-dithiophosphate | O,O-Dimethyl S-(2-(N-methylamino)-2-oxoethyl) dithiophosphate, phosphamide, Rogor | 0.03 | org. odr. | 4 |
| O,O-Dimethyl-O-(4-cyanphenyl)thiophosphate | Cyanox | 0.05 | org. odr. | 4 |
| O,O-Dimethyl-O-(3-methyl-4-nitrophenyl) thiophosphate | Methylnitrophos | 0.25 | org. odr. | 3 |
| O,O-Dimethyl-S- 2-(1-N-methylcarbomoyl-ethylmercapto) ethylthiophosphate | Kilval, vamidothion | 0.3 | org. odr. | 4 |
| N-(beta,beta-O,O-Diisopropyldithiophosphorylethyl) benzensulphonamide | O,O-Diisopropyl-S-2-phenylsulphonylaminoethylthiophosphate, Prefar, | 1.0 | s.-t. | 2 |

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| | Bensulide, Betasan | | | |
| 6.2.4. Salts of phosphoric acid and organic bases | | | | |
| 1,2,4-Triaminobenzene phosphate | | 0.01 | org. flavr. | 3 |
| n-aminobenzoic acid phosphate | | 0.1 | org. odr. | 3 |
| 7. Heterocyclic compounds | | | | |
| 7.1. Oxygen containing compounds | | | | |
| 7.1.1. Containing a three-partite cycle | | | | |
| Propylene oxyde | 1,2-Epoxy-propane, Methoxyrane | 0.01 | s.-t. | 2 |
| Epichlorhydrin | 1-Chlor--2,3 epoxypropane | 0.01 | s.-t. | 2 |
| 7.1.2. Containing a pentatomic cycle | | | | |
| Dichloromaleinic anhydride | Dichloro-butan-dionic anhydride | 0.1 | s.-t. | 2 |
| Furfurane | | 0.2 | s.-t. | 2 |
| 2-Methylfuran | Silvan | 0.5 | org. odr. | 4 |
| Furfuryl alcohol | Fur-2-ylmethanol, 2-hydroxymethylfuran, 2-furan-methanol | 0.6 <1> | s.-t. | 2 |
| Furfurol | 2-Furaldehyde | 1.0 | org. op. | 4 |
| 5-Nitrofurfurol-diacetate | (5-Nitro-2-furanyl) methanediol diacetate | 2.0 <1> | s.-t. | 2 |
| 7.1.3. Containing a hexatomic cycle | | | | |
| 5,6-Dihydro-4-methyl-2H-pyran | Methyldihydropyran | 0.0001 | s.-t. | 1 |

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| 4-Methyl-4-hydroxytetrahydro-pyran | 4-Methyltetrahydro-4-ol-2H-pyran, pyranic alcohol | 0.001 | s.-t. | 2 |
| Dimethyldioxane | 5,5-Dimethyl 1,3-dioxane | 0.005 | s.-t. | 2 |
| 4-Methyl-4-hydroxyethyl-1,3-dioxane | 4-Methyl-4-ethanol-1,3-dioxane, dioxanic alcohol | 0.04 | s.-t. | 2 |
| 7.1.4. Multinucleate | | | | |
| Chlorendic anhydride | Perchlorborn-5-ene-2,3-dicarbonic acid, anhydride | 1.0 | org. odr. | 3 |
| 7.2. Nitrogen containing compounds | | | | |
| 7.2.1. Pentatomic cycle with one nitrogen atom | | | | |
| Cyclohexylide of dichloromaleinic acid | Cimid | 0.04 | org. odr. | 4 |
| 7.2.2. Hexatomic aliphatic cycle with one nitrogen atom | | | | |
| Piperidine | | 0.06 | s.-t. | 3 |
| 4-Amino-2,2,6,6-tetramethylpiperidine | Triacetona-mine amine | 4.0 | s.-t. | 2 |
| Triacetona-mine | 2,2,6,6-Tetramethylpiperidin-4-one | 4.0 | s.-t. | 2 |
| 7.2.3. Hexatomic aliphatic cycle with one nitrogen atom | | | | |
| N-Methylpyridinium chloride | 1-Methylpyridinium chloride | 0.01 | org. odr. | 4 |
| Heptachloropicoline | 2-Trichloromethyl-3,4,5,6- | 0.02 | s.-t. | 2 |

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| | tetrachloro- pyridine | | | |
| Hexachloropicoline | 2-Trichloro- methyl-3,4, 5-trichloro- pyridine | 0.02 | s.-t. | 2 |
| Hexachloramino- picoline | 4-Amino-2- trichloro- methyl-3,5, 6-trichloro- pyridine | 0.02 | s.-t. | 2 |
| Pantachloramino- picoline | 4-Amino-2- trichloro- methyl-3,5, 6-dichloro- pyridine | 0.02 | s.-t. | 2 |
| Pentachloropicoline | 2-TRichloro- methyldichlo ropyridine | 0.02 | s.-t. | 2 |
| Tetrachloropicoline | 1-Chlor-6- (trichloro- methyl) pyridine | 0.02 | s.-t. | 3 |
| 2,5-Lutidine | 2,5-Dimethyl pyridine | 0.05 | s.-t. | 2 |
| alpha-Picoline | 2-Methyl- pyridine | 0.05 | s.-t. | 2 |
| Pyridine | | 0.2 | s.-t. | 2 |
| 4-Amino-3,5,6- trichlorpicolinic acid | 4-Amino- 3,5,6-tri- chlor-2- pyridincar- bonic acid, picloram, Tordon | 10.0 | s.-t. | 3 |
| Potassium 4-amino-3, 5,6-trichlor- picolinate | 4-Amino- 3,5,6-tri- chlor-2- pyridincar- bonic acid, potassium salt; Hloramp | 10.0 | s.-t. | 2 |
| 7.2.4. Multinucleate compounds with one nitrogen atom | | | | |

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| 5-Acetoxy-1,2-dimethyl-3-carbethoxyindole | Acetoxy-indole | 0.004 <1> | s.-t. | 2 |
| 6-Brom-5-hydroxy-3-carbethoxy-1-methyl-2-phenylthio-methylindole | Thioindole | 0.004 <1> | s.-t. | 2 |
| 2-Chlorcyclohexylthio-M-phthalimide | Phthalic acid, N-(2-chlorcyclohexylimide) | 0.02 | org. odr. | 4 |
| N-Trichloromethylthiophthalimide | Phthalan | 0.04 | org. odr. | 4 |
| 6-Brom-5-hydroxy-4-dimethylamino--carbethoxy-1-methyl-2-phenylthiomethylindole hydrochloride | Arbidol | 0.04 <1> | s.-t. | 3 |
| O,O-Dimethyl-S-phthalimidomethyl-dithiophosphate | Phthalophos | 0.2 | org. flavr. | 3 |
| Trichlormethylthio-tetrahydrophthalimide | Captan | 2.0 | org. odr. | 4 |
| 7.2.5. Pentatomic cycle with several nitrogen atoms | | | | |
| 1.3-Dichloro-5,5-dimethyl hydantoin | 5,5-Dimethyl 1,3-dichlorimidazolidin-2.4-dione, dichlorantane | None | s.-t. | 3 |
| 1-(2-Hydroxypropyl)-1-methyl-2-pentadecyl-2-imidazo-2-imidazolinium methyl sulphate | Carbosoline, SPD-3 | 0.2 | s.-t. | 2 |
| 1-Phenyl-3-pyrazolidone | Phenidone | 0.5 | org. clr. | 3 |
| 5,5-Dimethylhydantoin | | 1.0 | org. | 3 |

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| | | | flavr. | |
| 7.2.6. Pentatomic cycle with two nitrogen atoms | | | | |
| Sulphapyridazine | 6-(n-Amino-benzensulphamido)-3-methoxy-pyridazine; sul-sulphanilic acid, N-(6-methoxy-pyridazin-3-yl)amide | 0.2 <1> | s.-t. | 2 |
| O,O-Diethyl-O-(2-isopropyl-4-methylpyrimedyl-6-thiophosphate | O-(2-Isopropyl-6-methylpyrimidin-4-yl)-O,O-diethylthiophosphate, Bazudin | 0.03 | org. odr. | 4 |
| N-(2-Aminoethyl)piperazine | 1-(2-Aminoethyl)piperazine | 0.6 | s.-t. | |
| 1-Phenyl-4,5-dichlorpyridazone-6 | | 2.0 | s.-t. | 3 |
| 1-Phenyl-4-amino-5-chlorpyridazone-6 | 5-Amino-2-phenyl-4-chlorpyridazin-3(2H)-one, Phenazone | 2.0 | s.-t. | 2 |
| 4-Amino-6-chloropyrimidine | 6-Chlor-4-pyrimidinamine | 3.0 <1> | org. clr. | 3 |
| 4-Amino-6-methoxypyrimidine | | 5.0 <1> | org. clr. | 3 |
| Oxyethylpiperazine | | 6.0 | s.-t. | 2 |
| Diethylendiamine | Hexahydro-pyrazine, piperazine | 9.0 | org. odr. | 3 |
| 7.2.7. Pentatomic cycle with three nitrogen atoms | | | | |
| 2-Chlor-4,6-bis(ethylamino)-sym-triazine | 2,4-Bis(N-ethylamino)-6-chlor-1,3,5-tria- | None | org. flot. | 4 |

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| | zine, simazine | | | |
| 2-Chlor-4,6-bis(ethylamino)-sym-triazine 2-oxy derivative | 2-Oxy derivative of simazine | None | org. flot. | 4 |
| O,O-Dimethyl-S-(4,6-diamino-1,3,5-triazin-2-yl-methyl)-dithiophosphate | Sayfos, menazon, saphicol, azadition | 0.1 | s.-t. | 3 |
| Cyclotrimethylen-trinitroamine | 1,3,5-Tri-nitro-1,3,5-perhydrotri-azine, hexogene | 0.1 | s.-t. | 2 |
| 4,6-bis(Isopropylami-no)-2-(N-methyl-N-cyanamino)-1,3,5-triazine | Methazine | 0.3 | org. flavr. | 4 |
| 2-Amino-4-methyl-6-methoxy-1,3,5-triazine | 2-Amino-4-methyl-6-me-thoxy-sym-triazine | 0.4 <1> | org. odr. | 3 |
| 2-Chlor-4,6-bis(isopropylamino)-sym-triazine | 2,4-Bis(N-isopropyl-amino)-6-chlor-1,3,5-triazine, propazine, insoluble simazine | 1.0 | org. odr. | 4 |
| 2-Methylthio-4,6-diisopropylamino-sym-triazine | 2-Amino-4-(N,N-diiso-propylamino)-6-methyl-thio-1,3,5-tria-zine, pro-metryn | 3.0 | org. odr. | 3 |
| Cyanuric acid | 1,3,5-Tria-zine-2,4,6(1H,3H,5H)-trione | 6.0 | org. flavr. | 3 |
| 7.2.8. Multinucleate compounds with several nitrogen atoms | | | | |
| 1,2-Bis(1,4,6,9- | DKhTI 150A | 0.015 | s.-t. | 2 |

| | | | | |
|---|---|------|-----------|---|
| tetraazotricyclo-[4,4,1,1,4,9]-do-decano)-ethyliden dihydrochloride | | | | |
| Dipyridyl | Bipyridyl | 0.03 | org. odr. | 3 |
| 1,2,3-Benzotriazole | | 0.1 | s.-t. | 3 |
| Methyl-N-(2-benzimidazolyl) carbamate | 1H-benzimidazol-2-yl carbamic acid, methyl ether | 0.1 | org. film | 4 |
| 3-Cyclohexyl-5,6-trimethylenuracyl | 3-Cyclohexyl-6,7-dihydro-1H-cyclopentapyrimidin-2,4(3H,5H)-dione, Hexilur | 0.2 | s.-t. | 2 |
| 1,1-Dimethyl-4,4'-dipyridyldimethylphosphate | | 0.3 | org. odr. | 3 |
| Dipyridylphosphate | | 0.3 | org. odr. | 4 |
| Methyl-1-butylcarbamoyl-2-benzimidazolcarbamate | Arylate | 0.5 | org. film | 4 |
| Hexamethylenetetramine | 1,3,5,7-Tetraazatricyclo-decan, urotropine, amynoform, Formin | 0.5 | s.-t. | 2 |
| 5-Amino-2-(n-aminophenyl)-1H-benzimidazole | | 1.0 | s.-t. | 2 |
| Triethylendiamine | 1,4-Diazobicyclo[2.2.2]octane, DAVSO | 6.0 | s.-t. | 2 |
| 7.2.9. Containing over six atoms in the cycle | | | | |
| S-Ethyl-N-hexamethylenethio- | Hexahydro-1H-azepin- | 0.07 | org. odr. | 4 |

| | | | | |
|---|---|-------|----------------|---|
| carbamate | 1-thiocarbo- nic acid, S-ethyl ether; Yalan | | | |
| Hexamethylenamine hydrochloride | | 5.0 | s.-t. | 2 |
| Cyclotetramethylen- tetranitroamine | Octohydro- 1,3,5,7-tet- ranitro- 1,3,5,7-tet- razocyne, octogene | 0.2 | s.-t. | 2 |
| 7.3. Sulphur-containing compounds | | | | |
| 2-Chlorothiophene | | 0.001 | org. odr. | 4 |
| Tetrahydrothiophen- 1,1-dioxyde | Sulpholane, tetra- methylene sulphone | 0.5 | org. odr. | 3 |
| Thiophene | Thiofuran | 2.0 | org. odr. | 3 |
| 7.4. Mixed-type compounds | | | | |
| 7.4.1. Containing nitrogen and oxygen as heteroatoms | | | | |
| Codeine | | None | | |
| Morphine | | None | | |
| O,O-Diethyl-S-(6- chlorbenzoxazoliny- methyl)dithio- phosphate | S-(2,3-Di- hydro-3-oxo- -6-chlor- benzoxa- zol-3-ylme- thyl)-O,O- diethylphos- phate, Phosalone | 0.001 | org. odr. | 4 |
| Tetrahydro-1,4- oxazine | Morpholine | 0.04 | org. flavr. | 3 |
| Benzoxazolone-2 | benzoxazol- 2(3H)-one | 0.1 | s.-t. | 2 |
| 3-Chlormetal-6- chlorbenzoxazolone | 6-Chlor-3- chlormethyl -2-(3H)ben- zoxazolone | 0.4 | s.-t. | 2 |

| 7.4.2. Containing nitrogen and sulphur as heteroatoms | | | | |
|---|---|----------|-----------|---|
| Dibenzthiazoldi-sulphide | 2,2'-Dithio-dibenzothiazole, altax | None | org. odr. | 3 |
| 2-Butylthiobenzothiazole | Butylcaptax | 0.005 | org. odr. | 4 |
| 3,5-Dimethyltetrahydro-1,3,5-thiadiazinethione-2 | 3,5-Dimethylperhydro-1,3,5-thiadiazin-2-thione, mylon, thiazone | 0.01 | org. odr. | 4 |
| Benzothiazole | | 0.25 <1> | org. odr. | 4 |
| 2-Hydroxybenzothiazole | 2-(3H)-Hydroxybenzothiazolone | 1.0 | s.-t. | 2 |
| 2-Mercaptobenzothiazole | Benzothiazol-2-thiol, Captax | 5.0 | org. odr. | 4 |
| 8. Organoelemental compounds | | | | |
| 8.1. Mercury compounds | | | | |
| Ethyl mercuric chloride | Granosan | 0.0001 | s.-t. | 1 |
| Mercury diethyl | | 0.0001 | s.-t. | 1 |
| 8.2. Tin compounds | | | | |
| Tetraethyltin | Tetraethylstannane | 0.0002 | s.-t. | 1 |
| Bis(tributyltin) oxyde | | 0.0002 | s.-t. | 1 |
| Tin tributylmethacrylate | Tributyl(2-methyl-1-oxo-2-propenyl) oxystannane | 0.0002 | s.-t. | 1 |
| Dicyclohexyltin oxyde | Dicyclohexyloxostannane | 0.001 | s.-t. | 2 |
| Tricyclohexyltin chloride | | 0.001 | s.-t. | 2 |

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| Dichlordibutyltin | Dibutyldi-chlor-stannane | 0.002 | s.-t. | 2 |
| Diethyltin dichloride | Dichlor-diethyl-stannane | 0.002 | s.-t. | 2 |
| Tetrabutyltin | Tetrabutyl-stannane | 0.002 | s.-t. | 2 |
| Ethylene-bis(thioglycolate)-dioctyltin | | 0.002 | s.-t. | 2 |
| Dibutyltin oxyde | Dibutyloxo-stannane | 0.004 | s.-t. | 2 |
| Dibutyltin dilaurate | Bis(dodecanoyloxy)-dinbutyl-stannane | 0.01 | s.-t. | 2 |
| Dibutyltin diisooctylthioglycolate | Bis(isooctylloxycarbonylmethylthio)dibutylstannane | 0.01 | s.-t. | 2 |
| Diethyltin dioctanoate | Diethyl-bis(octanoyloxy)stannane, tin diethyl-dicaprylate | 0.01 | s.-t. | 2 |
| Diisobutyltin maleatedioctyl | | 0.02 | s.-t. | 2 |
| Sulphidedibutyltin | Dibutyltin sulphide | 0.02 | s.-t. | 2 |
| Tributyltin chloride | Chlortributylstannane, tributyl-chlor-stannane | 0.02 | s.-t. | 2 |
| 8.3. Lead compounds | | | | |
| Tetraethyllead | | None | s.-t. | 1 |
| 8.4. Arsenic compounds | | | | |
| 8.5. Silicium compounds | | | | |

| | | | |
|-----------------------|-----|----------------|---|
| Trifluoropropylsilane | 1.5 | org. flavr. | 4 |
|-----------------------|-----|----------------|---|

Supplement 3
(for reference)

ALPHABETICAL INDEX OF HARMFUL SUBSTANCES
IN DRINKING WATER SPECIFIED IN SUPPLEMENT 2

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| Acrylic aldehyde | 3.2.1.1.2. |
| Acricide | 4.2.1.2.2.1.2. |
| Acrolein | 3.2.1.1.2. |
| Alizarin | 3.2.2.2. |
| Alkylaminopropionitrile C17 - C20 | 4.1.3.1.1. |
| Alkylaniline | 4.1.1.1.2.2.1. |
| Alkylbenzyldimethylammonium chloride C10 - C16 | 4.1.3.2.2. |
| Alkylbenzyldimethylammonium chloride C17 - C20 | 4.1.3.2.2. |
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| Alkylpropylendiamine | 4.1.1.2.1.2. |
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| N-(C7-C9)Alkyl-N-phenyl-n-phenylendiamine | 4.1.3.2.2. |
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| Aldrin | 2.2.1.2. |
| Altax | 7.4.2. |
| Amylcarbinol | 3.1.1.1. |
| 5-Amino-2-(n-aminophenyl)-1H-benzimidazole | 7.2.8. |
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| Aminobenzene | 4.1.1.1.2.2.1. |
| 3-Aminobenzotrifluoride | 4.1.1.1.2.2.1.1. |
| n-Aminobutylbenzene | 4.1.1.1.2.2.1. |
| 6-(n-Aminobenzensulphamido)-3-methoxy- pyridazine | 7.2.6. |

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| 1-Amino-2-hydroxybenzene | 4.1.1.1.2.2.1.2. |
| 1-Amino-3-hydroxybenzene | 4.1.1.1.2.2.1.2. |
| 1-Amino-2-hydroxypropane | 4.1.1.1.2.2.1.2. |
| 4-Amino-2-(2-hydroxyethyl)-N-ethylanilin-sulphite | 4.1.2.2.2.1. |
| 2-Amino-4-(N,N-diisopropylamino)-6-methylthio-1,3,5-triazine | 7.2.7. |
| 4-Aminodiphenylamine | 4.1.2.2.2. |
| 1,4-Aminodiethylaniline sulphate | 4.1.3.2.2. |
| n-Aminomethylbenzene | 4.1.1.1.2.2.1. |
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| Potassium 4-amino-3,5,6-trichloropicolinate | 7.2.3. |
| Aminophenetol | 4.1.1.1.2.2.1.2. |
| 5-Amino-2-phenyl-4-chlorpyridazin-3(2H)one | 7.2.6. |
| m-Aminophenol | 4.1.1.1.2.2.1.2. |
| o-Aminophenol | 4.1.1.1.2.2.1.2. |
| n-Aminophenol | 4.1.1.1.2.2.1.2. |
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| n-Anisidine | 4.1.1.1.2.2.1.2. |
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| 5-(2-Acetamidoethyl)-O,O-dimethyldithiophosphate | 6.2.2.2. |
| n-Acetaminophenol | 4.1.2.2.2.1. |
| N-Acetyl-2-aminophenol | 4.1.2.2.2.1. |

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| Acetylacetonates | 3.2.2. |
| 5-Acetoxy-1,2-dimethyl-3-carbethoxyindole | 7.2.4. |
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| -B- | |
| Bazudin | 7.2.6. |
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| Benzamide | 4.1.1.1.2.2.1.3. |
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| 5-Benzyl-O,0-diethylthiophosphate | 6.2.2.2. |
| 3-Benzyltoluene | 1.2.2.1. |
| Benzyl chloride | 2.2.2.1.2. |
| Benzyl cyanide | 4.1.3.2.2.1. |
| Benzoxazol-2(3H)-one | 7.4.1. |
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| Benzene | 1.2.2.1. |
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| Benzensulphamide | 5.2.4.1.4. |
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| 2,2-Bis-(4-hydroxy-3,5-dichlorophenyl)propane | 3.1.3.2.1. |
| Bis(2-hydroxyethyl)methylamine | 4.1.3.1.2. |
| Bis(dodecanouloxy)-di-n-butylstannane | 8.2. |
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| 4,6-Bis(isopropylamino)-2-(N-methyl-N-cyanamino)-1,3,5-triazine | 7.2.7. |
| 2,4-Bis(N-isopropylamino)-6-chlor-1,3,5-triazine | 7.2.7. |
| Bis(2-methylpropyl)amine | 4.1.2.1. |
| 1,4-Bis(4-methyl-2-sulphophenylamino)-5,8-dihydroxyanthraquinone, disodium salt | 5.2.4.1.1.1. |
| 1,2-Bis-methoxycarbonyl thioureidobenzene | 5.1.4.1. |
| 1,2-Bis(1,4,6,9-tetraazotricyclo[4,4,1,1,-4,9]dodecano)-ethyliden dihydrochloride | 7.2.8. |
| Bis(tributyltin)oxyde | 8.2. |

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| 1,3-Bis(trichlormethyl)benzene | 2.2.2.1.2. |
| 1,4-Bis(trichlormethyl)benzene | 2.2.2.1.2. |
| Bis(n-chlorophenyl)sulphone | 5.2.2. |
| O,O-Bis(2-chlorethyl)vinylphosphonate | 6.1.3. |
| 2,4-Bis(N-ethylamino)-6-chlor-1,3,5-triazine | 7.2.7. |
| Bicyclo(2,2,1)hepta2,5-diene | 1.2.1.2. |
| Biphenyl | 1.2.2.2.1. |
| Bolstar | 6.2.2.2. |
| Botran | 4.2.1.2.2.1.3.1. |
| 3-Brombenzaldehyde | 3.2.1.2.2.1.1. |
| m-Brombenzaldehyde | 3.2.1.2.2.1.1. |
| 6-Brom-5-hydroxy-4-dimethylamino-3-carb-ethoxy-1-methyl-2-phenylthiomethylindole hydrochloride | 7.2.4. |
| 6-Brom-5-hydroxy-3-carbethoxy-1-methyl-2-phenyl-thiomethylindole | 7.2.4. |
| O-(4-Brom-2,5-dichlorophenyl)-O,O-dimethyl thiophosphate | 6.2.2.2.1. |
| Bromcamphora | 3.2.1.2.1.1. |
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| Bromophos | 6.2.2.2.1. |
| Ethyl bromide | 2.1.1. |
| Bromtoluidine(mix of o,m,n-isomers) | 4.1.1.1.2.2.1.1. |
| Bromotoluene | 4.1.1.1.2.2.1.1. |
| Butadiene-1,3 | 1.1. |
| Butamide | 5.2.2. |
| Butandinitrile | 4.1.3.1.1. |
| 1,4-Butandiol | 3.1.3.1. |
| Butan-1,4-diol | 3.1.3.1. |
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| Butan-2-ol | 3.1.1.1. |
| Butan-2-one | 3.2.1.1.1. |
| But-1-ene | 1.1. |
| 2-Butenal | 3.2.1.1.2. |
| But-2-enal | 3.2.1.1.2. |
| Butyl acrylate | 3.3.2.1.1.2.1. |
| n-Butylamide of benzenesulphonic acid | 5.2.4.1.4. |
| Butylamide of O-ethyl-S-phenyldithio-phosphoric acid | 6.2.2.2.2. |
| Butylamine | 4.1.1.1.1.1. |
| t-Butylamine | 4.1.1.1.1.1. |
| n-Butylaniline | 4.1.1.1.2.2.1. |
| Butylbenzene | 1.2.2.1. |
| N-Butylbenzensulphamide | 5.2.4.1.4. |
| 1-Butylbiguanidine hydrochloride | 4.1.3.1. |
| N-Butyl-1-butanamine | 4.1.2.1. |
| 2-sec-Butyl-4,6-dinitrophenyl-3,3-dimethyl acrylate | 4.2.1.2.2.1.2. |
| 2-sec-Butyl-4,6-dinitrophenyl-3-methylcro-tonate | 4.2.1.2.2.1.2. |
| Butylene | 1.1. |
| Butylcarbinol | 3.1.1.1. |
| Butylcaptax | 7.4.2. |
| Butylxanthogenate | 5.1.4.3. |

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| N-n-Butyl-N-(n-methylbenzensulphonyl)-urea | 5.2.2. |
| Butylnitrite | 4.2.2. |
| t-Butyl alcohol | 3.1.1.1. |
| Butyl ether of acrylic acid | 3.3.2.1.1.2.1. |
| Butyl ether 2,4-D | 3.3.2.1.1.1.3.1. |
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| Vinylcarbinol | 3.1.1.1. |
| Vinyl ether of monoethanolamine | 4.1.1.1.1.2.1. |
| Vinyl ether of acetic acid | 3.3.2.1.1.1.1.2. |
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| Hexahydrobenzene | 1.2.1.1. |
| 1,4,4a,5,8,8a-Hexahydro -1,2,3,4,10,10 -hexachlor-1,4,5,8-dimethanenaphthalene | 2.2.1.2. |
| 2,3,3a,4,7,7a-Hexahydro -2,4,5,6,7,8,8 -heptachlor-4,7-methanindene | 2.2.1.2. |
| 3-(Hexahydro -4,7-methanindan-5-yl)-1,1 -dimethylurea | 4.1.3.2.1. |
| Hexahydropyrazine | 7.2.6. |
| Hexahydrophenol | 3.1.1.2.1. |
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| Hexamethylenediamine | 4.1.1.2.1.1. |
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| Hexachlorbutadiene | 2.1.2. |
| 1,2,3,4,10,10-Hexachlor-1,4,4a,5,8,8a-he- | 2.2.1.2. |

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| xahydro-1,4-endo-exo-5,8-dimethanenaphthalene | |
| Hexachlorometaxylene | 2.2.2.1.2. |
| Hexachloroparaxylene | 2.2.2.1.2. |
| Hexachloropicoline | 7.2.3. |
| 1,2,3,4,5,6-Hexachlorocyclohexane | 2.2.1.1. |
| Hexachlorocyclopentadiene | 2.2.1.1. |
| 1,2,3,4,5,5-Hexachloro-1,3-cyclopentadiene | 2.2.1.1. |
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| Hexogene | 7.2.7. |
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