

Растворители для GC-MS, безводные, дейтерированные для NMR

Технические характеристики

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SOLVENTS FOR GC-MS



Solvents for GC-MS are crucial chemical substances in gas chromatography-mass spectrometry. They are used to dissolve samples before injection into the system, facilitating the separation and analysis of components. These solvents, must be of high purity to avoid interference with analyte detection. The choice of solvent depends on the nature of the sample and specific analytical requirements.



2,2,4-Trimethylpentane, GC-MS

- Synonyms: Isooctane, Isobutyltrimethylmethane, iso-Octane
- C_8H_{18}
- $M = 114,26 \text{ g/mol}$
- CAS [540-84-1]
- EINECS-No.: 208-759-1
- Density: $0,69 \text{ g/cm}^3$
- Solub. in water: (25 °C): $0,56 \text{ mg/l}$
- Melting point: -107 °C
- Boiling point: 99 °C
- Flash pt. -12 °C
- Ignition temp.: 410 °C
- Vapour pressure: (20 °C) 51 hPa
- Dielectric const.: (20 °C) $1,9$
- LD 50 (oral, rat): $> 2000 \text{ mg/kg}$
- EC-Index-No.: 601-009-00-8
- ADR: 3 F1 II UN 1262
- IMDG: 3 II UN 1262
- IATA/ICAO: 3 II UN 1262
- GHS-signal word: Danger
- GHS-H sentences: H225 - H304 - H400 - H410 - H315 - H336
- GHS-P sentences: P210 - P301+P310 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2901 10 00 00
- Applications: analytical chemistry, solvent for fat and oil extractions; in determining octane numbers of fuels.

SPECIFICATIONS

assay (G.C.): min. 99,5 %

colour (Hazen): max. 10

identity (IR-spectrum): passes test

residue on evaporation: max. 3 ppm
water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis

Volume

x 1 l

Reference

[IS01671000](#)

Packaging

x 1 l :: Glass bottle

Volume

x 2,5 l

Reference

[IS01672500](#)

Packaging

x 2,5 l :: Glass bottle



Acetone, GC-MS

- Synonyms: Dimethyl ketone, 2-Propanone
- C₃H₆O
- M = 58,08 g/mol
- CAS [67-64-1]
- EINECS-No.: 200-662-2
- Density: 0,79 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: -95 °C
- Boiling point: 56 °C
- Flash pt. < -20 °C
- Ignition temp.: 540 °C
- Vapour pressure: (20 °C) 233 hPa
- Refraction index: (n 20 °C/D) 1,3588
- Dielectric const.: (25 °C) 20,7
- LD 50 (oral, rat): 5800 mg/kg
- EC-Index-No.: 606-001-00-8
- ADR: 3 F1 II UN 1090
- IMDG: 3 II UN 1090
- IATA/ICAO: 3 II UN 1090
- GHS-signal word: Danger
- GHS-H sentences: H225 - H319 - H336 - EUH066
- GHS-P sentences: P210 - P303+P361+P353 - P305+P351+P338 - P370+P378 - P405 - P501a
- Tariff number: 2914 11 00 00
- Applications: solvents, analytical chemistry, synthesis of organic products, photography.

SPECIFICATIONS

assay (G.C.): min. 99,8 %

colour (Hazen): max. 10

identity (IR-spectrum): passes test

residue on evaporation: max. 3 ppm

water (K.F.) : max. 0,2 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual

signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Acetonitrile, GC-MS

- Synonyms: Methyl cyanide, Cyanomethane
- CH3CN
- M = 41,05 g/mol
- CAS [75-05-8]
- EINECS-No.: 200-835-2
- Density: 0,786 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: -45,7 °C
- Boiling point: 81,6 °C
- Flash pt. 2 °C
- Ignition temp.: 524 °C
- Vapour pressure: (20 °C) 97 hPa
- Refraction index: (n 20 °C) 1,3442
- Dielectric const.: (20 °C) 37,5
- LD 50 (oral, rat): 2730 - 3800 mg/kg
- EC-Index-No.: 608-001-00-3
- ADR: 3 F1 II UN 1648
- IMDG: 3 II UN 1648
- IATA/ICAO: 3 II UN 1648
- GHS-signal word: Danger
- GHS-H sentences: H225 - H302+H312+H332 - H319
- GHS-P sentences: P210 - P303+P361+P353 - P305+P351+P338 - P370+P378 - P403+P235
- P501a
- Tariff number: 2926 90 70 22
- Applications: chromatography, synthesis of organic products, solvents.

SPECIFICATIONS

assay (G.C.): min. 99,8 %

colour (Hazen): max. 10

identity (IR-spectrum): passes test

residue on evaporation: max. 3 ppm

water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Cyclohexane, GC-MS

- Synonyms: Hexahydrobenzene, Hexamethylene, Naphthene
- C6H12
- M = 84,16 g/mol
- CAS [110-82-7]
- EINECS-No.: 203-806-2
- Density: 0,78 g/cm³
- Solub. in water: (20 °C): 55 mg/l
- Melting point: 6 °C
- Boiling point: 80,7 - 81 °C
- Flash pt. -18 °C

- Ignition temp.: 260 °C
- Vapour pressure: (20 °C) 103 hPa
- Refraction index: (n 20 °C/D) 1,4264
- Dielectric const.: (20 °C) 2,0
- LD 50 (oral, rat): 12705 mg/kg
- EC-Index-No.: 601-017-00-1
- ADR: 3 F1 II UN 1145
- IMDG: 3 II UN 1145
- IATA/ICAO: 3 II UN 1145
- GHS-signal word: Danger
- GHS-H sentences: H225 - H315 - H319 - H335 - H336 - H304 - H410
- GHS-P sentences: P210 - P301+P310 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2902 11 00 00
- Applications: solvents, analytical chemistry, synthesis of organic products.

SPECIFICATIONS

assay (G.C.): min. 99,8 %
 colour (Hazen): max. 10
 identity (IR-spectrum): passes test
 residue on evaporation: max. 3 ppm
 water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Dichloromethane, GC-MS, suitable for nitrosamine analysis

- Synonyms: Methylene chloride, Chloromethylene
- CH2Cl2
- M = 84,93 g/mol
- CAS [75-09-2]
- EINECS-No.: 200-838-9
- Density: 1,32 g/cm³
- Solub. in water: (20 °C): 20 g/l
- Melting point: ~ -95 °C
- Boiling point: 40 °C
- Ignition temp.: 605 °C
- Vapour pressure: (20 °C) 475 hPa
- Dielectric const.: (20 °C) 9,1
- LD 50 (oral, rat): 1600 mg/kg
- EC-Index-No.: 602-004-00-3
- ADR: 6.1 T1 III UN 1593
- IMDG: 6.1 III UN 1593
- IATA/ICAO: 6.1 III UN 1593
- GHS-signal word: Warning
- GHS-H sentences: H315 - H319 - H335 - H336 - H351 - H373 -
- GHS-P sentences: P260 - P280 - P305+P351+P338 - P321 - P405 - P501a
- Tariff number: 2903 12 00 00
- Applications: solvents, analytical chemistry.

SPECIFICATIONS

assay (G.C.): min. 99,8 %
 colour (Hazen): max. 10
 identity (IR-spectrum): passes test
 nitrosamine analysis: passes test
 residue on evaporation: max. 3 ppm
 water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Ethyl acetate, GC-MS

- Synonyms: Acetic acid ethyl ester, Acetic ether
- $C_4H_8O_2$
- $M = 88,10 \text{ g/mol}$
- CAS [141-78-6]
- EINECS-No.: 205-500-4
- Density: $0,90 \text{ g/cm}^3$
- Solub. in water: (20 °C): $85,3 \text{ g/l}$
- Melting point: -83 °C
- Boiling point: 77 °C
- Flash pt. -4 °C
- Ignition temp.: 460 °C
- Vapour pressure: (20 °C) 97 hPa
- Refraction index: (n 20 °C/D) $1,3723$
- Dielectric const.: (25 °C) $6,0$
- LD 50 (oral, rat): 5620 mg/kg
- EC-Index-No.: 607-022-00-5
- ADR: 3 F1 II UN 1173
- IMDG: 3 II UN 1173
- IATA/ICAO: 3 II UN 1173
- GHS-signal word: Danger
- GHS-H sentences: H225 - H319 - H336 - EUH066
- GHS-P sentences: P210 - P303+P361+P353 - P305+P351+P338 - P370+P378 - P405 - P501a
- Tariff number: 2915 31 00 00
- Applications: solvents, perfumery, photography.

SPECIFICATIONS

assay (G.C.): min. 99,8 %
colour (Hazen): max. 10
identity (IR-spectrum): passes test
residue on evaporation: max. 3 ppm
water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Methanol, GC-MS

- Synonyms: Methyl alcohol, Carbinol, Methynol, Wood alcohol
- CH_3OH
- $M = 32,04 \text{ g/mol}$
- CAS [67-56-1]
- EINECS-No.: 200-659-6
- Density: $0,792 \text{ g/cm}^3$
- Solub. in water: (20 °C): miscible
- Melting point: -98 °C

- Boiling point: 65 °C
- Flash pt. 10 °C
- Ignition temp.: 455 °C
- Vapour pressure: (20 °C) 128 hPa
- Refraction index: (n 20 °C/D) 1,3288
- Dielectric const.: (25 °C) 32,6
- LD 50 (oral, rat): 5628 mg/kg
- EC-Index-No.: 603-001-00-X
- ADR: 3 FT1 II UN 1230
- IMDG: 3 II UN 1230
- IATA/ICAO: 3 II UN 1230
- GHS-signal word: Danger
- GHS-H sentences: H225 - H301 - H311 - H331 - H370
- GHS-P sentences: P210 - P301+P310 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2905 11 00 10
- Applications: solvents, synthesis of organic products, in antifreeze compositions, solvent for animal and vegetable oils extractions.
- Appearance: Incolore

SPECIFICATIONS

assay (G.C.): min. 99,0 %
 colour (Hazen): max. 10
 identity (IR-spectrum): passes test
 residue on evaporation: max. 3 ppm
 water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



n-Hexane, GC-MS

- Synonyms: n-Caproylhydride, n-Hexylhydride
- C₆H₁₄
- M = 86,18 g/mol
- CAS [110-54-3]
- EINECS-No.: 203-777-6
- Density: 0,66 g/cm³
- Solub. in water: (20 °C): 0,0095 g/l
- Melting point: -94,3 °C
- Boiling point: 69 °C
- Flash pt. -22 °C
- Ignition temp.: 240 °C
- Vapour pressure: (20 °C) 160 hPa
- Dielectric const.: (20 °C) 1,8
- LD 50 (oral, rat): 28710 mg/kg
- EC-Index-No.: 601-037-00-0
- ADR: 3 F1 II UN 1208
- IMDG: 3 II UN 1208
- IATA/ICAO: 3 II UN 1208
- GHS-signal word: Danger
- GHS-H sentences: H225 - H304 - H361f - H373 - H315 - H336 - H411
- GHS-P sentences: P210 - P301+P310 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2901 10 00 00

SPECIFICATIONS

assay (G.C.): min. 96,0 %
 colour (Hazen): max. 10
 identity (IR-spectrum): passes test

residue on evaporation: max. 3 ppm
water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis



Toluene, GC-MS

- Synonyms: Methylbenzene, Phenylmethane
- C₇H₈
- M = 92,14 g/mol
- CAS [108-88-3]
- EINECS-No.: 203-625-9
- Density: 0,87 g/cm³
- Solub. in water: (20 °C): 0,52 g/l
- Melting point: -95 °C
- Boiling point: 111 °C
- Flash pt. 4 °C
- Ignition temp.: 535 °C
- Vapour pressure: (20 °C) 29 hPa
- Dielectric const.: (25 °C) 2,3
- LD 50 (oral, rat): 636 mg/kg
- EC-Index-No.: 601-021-00-3
- ADR: 3 F1 II UN 1294
- IMDG: 3 II UN 1294
- IATA/ICAO: 3 II UN 1294
- GHS-signal word: Danger
- GHS-H sentences: H225 - H304 - H361d - H373 - H315 - H336
- GHS-P sentences: P210 - P301+P310 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2902 30 00 00
- Applications: synthesis of organic products, solvents, as gasoline additive.

SPECIFICATIONS

assay (G.C.): min. 99,8 %
colour (Hazen): max. 10
identity (IR-spectrum): passes test
residue on evaporation: max. 3 ppm
water (K.F.) : max . 0,05 %

GC/MSD (retention range n-undecane to n-tetracontane, scanning area 30 - 600 amu, individual signals (n-tetradecane standard)): max. 3,0 ng/ml (ppb)

Suitable for residue analysis

SOLVENTS, ANHYDROUS



Anhydrous solvents are characterized by a very low water content. They are used in processes or syntheses that are sensitive to the presence of any water.



1,4-Dioxane, 99%, anhydrous (max. 0,005% H₂O), with molecular sieves, stabilized with 2,5 ppm of 2,6-Di-tert-butyl-4-methylphenol (BHT)



1,4-Dioxane, 99,5%, anhydrous (max. 0,005% H₂O), stabilized with 2,5 ppm of 2,6-Di-tert-butyl-4-methylphenol (BHT)



2,2,4-Trimethylpentane, 99,5%, anhydrous (max. 0,003% H₂O)



2-Propanol, 99,8%, anhydrous (max. 0,005% H₂O)



Acetone, 99,8%, anhydrous (max. 0,005% H₂O)



Acetonitrile, 99,7%, anhydrous (max. 0,005% H₂O), with molecular sieves



Acetonitrile, 99,9%, anhydrous (max. 0,001% H₂O)



Benzyl alcohol, 99,5%, anhydrous (max. 0,01% H₂O)



Chloroform, 99,9%, anhydrous (max. 0,003% H₂O), stabilized with 150 ppm of amylene



Chloroform, 99,9%, anhydrous (max. 0,003% H₂O), with molecular sieves, stabilized with 150 ppm of amylene



Dichloromethane, 99,9%, anhydrous (max. 0,003% H₂O), stabilized with approx. 50 ppm of amylene



Dichloromethane, 99,9%, anhydrous (max. 0,003% H₂O), with molecular sieves



Diethyl ether, 99,5%, anhydrous (max. 0,005% H₂O), with molecular sieves, stabilized with approx. 7 ppm of 2,6-Di-tert-butyl-4-methylphenol (BHT)



Diethyl ether, 99,7%, anhydrous (max. 0,005% H₂O), stabilized with approx. 7 ppm of 2,6-Di-tert-butyl-4-methylphenol (BHT)



Dimethyl sulfoxide, 99,5%, anhydrous (max. 0,005% H₂O), with molecular sieves



Dimethyl sulfoxide, 99,9%, anhydrous (max. 0,005% H₂O)



Ethyl acetate, 99,8%, anhydrous (max. 0,005% H₂O), with molecular sieves



Methanol, 99,9%, anhydrous (max. 0,003% H₂O)



N,N-Dimethylacetamide, 99,5%, anhydrous (max. 0,005% H₂O)



N,N-Dimethylformamide, 99,8%, anhydrous (max. 0,005% H₂O)



n-Hexane, 96%, anhydrous (max. 0,002% H₂O)



Pyridine, 99,5%, anhydrous (max. 0,005% H₂O)



Pyridine, 99,5%, anhydrous (max. 0,005% H₂O), with molecular sieves



Tetrahydrofuran, 99,5%, anhydrous (max. 0,005% H₂O), with molecular sieves, stabilized with 2,6-Di-tert-butyl-4-methylphenol (BHT)



Tetrahydrofuran, 99,8%, anhydrous (max. 0,005% H₂O), stabilized with 250 ppm of 2,6-Di-tert-butyl-4-methylphenol (BHT)



Toluene, 99,5%, anhydrous (max. 0,005% H₂O), with molecular sieves



Toluene, 99,8%, anhydrous (max. 0,003% H₂O)

SOLVENTS, DEUTERATED FOR NMR



In deuterated solvents, the hydrogen atoms are replaced with deuterium. Deuterium is a heavy isotope of hydrogen with a proton and a neutron in its nucleus, instead of a single proton as in hydrogen. Deuterated solvents produce a sharper signal in nuclear magnetic resonance (NMR) spectra. They are commonly used in NMR spectroscopy and other chemical and physical analytical techniques.



Acetone-d₆, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Hexadeuteroacetone
- C₃D₆O
- M = 64,12 g/mol
- CAS [666-52-4]
- EINECS-No.: 211-563-9
- Density: 0,87 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: -95,4 °C
- Boiling point: 56 °C
- Flash pt. < -20 °C
- Ignition temp.: 540 °C
- Vapour pressure: (20 °C) 233 hPa
- Dielectric const.: (25 °C) 20,7
- LD 50 (oral, rat): 5800 mg/kg
- EC-Index-No.: 606-001-00-8
- ADR: 3 F1 II UN 1090
- IMDG: 3 II UN 1090
- IATA/ICAO: 3 II UN 1090
- GHS-signal word: Danger
- GHS-H sentences: H225 - H319 - H336 - EUH066
- GHS-P sentences: P210 - P303+P361+P353 - P305+P351+P338 - P370+P378 - P405 - P501a
- Tariff number: 2845 90 10 00

- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,03 %

performance test (NMR-spectrum): passes test



Acetonitrile-d₃, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Trideuteroacetonitrile
- CD₃CN
- M = 44,05 g/mol
- CAS [2206-26-0]
- EINECS-No.: 218-616-5
- Density: 0,84 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: -46 °C
- Boiling point: 79 °C
- Flash pt. 5 °C
- Ignition temp.: 525 °C
- Vapour pressure: (20 °C) 97 hPa
- LD 50 (oral, rat): 2730 - 3800 mg/kg
- ADR: 3 F1 II UN 1648
- IMDG: 3 II UN 1648
- IATA/ICAO: 3 II UN 1648
- GHS-signal word: Danger
- GHS-H sentences: H225 - H302 - H312 - H332 - H319
- GHS-P sentences: P210 - P303+P361+P353 - P305+P351+P338 - P370+P378 - P403+P235 - P501a
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,03 %

performance test (NMR-spectrum): passes test



Chloroform-d + TMS (99:1, v/v), deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- CDCl₃
- M = 120,38 g/mol
- CAS [865-49-6]
- EINECS-No.: 212-742-4
- Density: 1,50 g/cm³
- Solub. in water: (20 °C): 8,2 g/l
- Melting point: -64,1 °C
- Boiling point: 60 °C
- Vapour pressure: (20 °C) 211 hPa
- LD 50 (oral, rat): 908 mg/kg
- EC-Index-No.: 602-006-00-4
- ADR: 6.1 T1 III UN 1888
- IMDG: 6.1 III UN 1888
- IATA/ICAO: 6.1 III UN 1888

- GHS-signal word: Danger
- GHS-H sentences: H331 - H372 - H351 - H361d - H302 - H315 - H319
- GHS-P sentences: P260 - P280 - P305+P351+P338 - P321 - P405 - P501a
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,02 %

performance test (NMR-spectrum): passes test



Chloroform-d, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- CDCl₃
- M = 120,38 g/mol
- CAS [865-49-6]
- EINECS-No.: 212-742-4
- Density: 1,50 g/cm³
- Solub. in water: (20 °C): 8,2 g/l
- Melting point: -64,1 °C
- Boiling point: 60 °C
- Vapour pressure: (20 °C) 211 hPa
- LD 50 (oral, rat): 908 mg/kg
- EC-Index-No.: 602-006-00-4
- ADR: 6.1 T1 III UN 1888
- IMDG: 6.1 III UN 1888
- IATA/ICAO: 6.1 III UN 1888
- GHS-signal word: Danger
- GHS-H sentences: H331 - H372 - H351 - H361d - H302 - H315 - H319
- GHS-P sentences: P260 - P280 - P305+P351+P338 - P321 - P405 - P501a
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,01 %

performance test (NMR-spectrum): passes test



Deuterium oxide, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Heavy water
- D₂O
- M = 20,03 g/mol
- CAS [7789-20-0]
- EINECS-No.: 232-148-9
- Density: 1,11 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: 3,8 °C
- Boiling point: 101 °C
- Tariff number: 2845 10 00 00
- Applications: for nuclear magnetic resonance spectroscopy, for nuclear reactions.

SPECIFICATIONS

deuteration degree: min. 99,8 %
performance test (NMR-spectrum): passes test



Dichloromethane-d₂, deuteration degree min. 99,5%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Methylene chloride-d₂, Dideuteromethylene chloride, Dideuterodichloromethane
- CCl₂D₂
- M = 86,95 g/mol
- CAS [1665-00-5]
- EINECS-No.: 216-776-0
- Density: 1,36 g/cm³
- Solub. in water: (20 °C): 20 g/l
- Melting point: -97 °C
- Boiling point: 39 °C
- Ignition temp.: 605 °C
- Vapour pressure: (20°C) ~ 450 hPa
- LD 50 (oral, rat): 1600 mg/kg
- EC-Index-No.: 602-004-00-3
- ADR: 6.1 T1 III UN 1593
- IMDG: 6.1 III UN 1593
- IATA/ICAO: 6.1 III UN 1593
- GHS-signal word: Warning
- GHS-H sentences: H351
- GHS-P sentences: P280 - P201 - P202 - P308+P313 - P405 - P501a
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,5 %
water (K.F., H₂O + D₂O): max. 0,03 %
performance test (NMR-spectrum): passes test



Dimethyl sulfoxide-d₆, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Methylsulfoxide deuterated, DMSO deuterated, Hexadeuterodimethyl sulfoxide
- C₂D₆OS
- M = 84,17 g/mol
- CAS [2206-27-1]
- EINECS-No.: 218-617-0
- Density: 1,19 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: 20,2 °C
- Boiling point: 190 °C
- Flash pt. 88 °C
- Ignition temp.: 270 °C
- Vapour pressure: (20 °C) 2,5 hPa
- Refraction index: (20 °C) 1,48
- LD 50 (oral, rat): 17500 mg/kg
- GHS-signal word: Warning
- GHS-H sentences: H315 - H319
- GHS-P sentences: P280 - P264 - P305+P351+P338 - P321 - P332+P313 - P337+P313
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,03 %

performance test (NMR-spectrum): passes test



Methanol-d₄, deuteration degree min. 99,8%, NMR spectroscopy grade, Spectrosol®

- Synonyms: Tetradeuteromethanol
- CD₃OD
- M = 36,07 g/mol
- CAS [811-98-3]
- EINECS-No.: 212-378-6
- Density: 0,89 g/cm³
- Solub. in water: (20 °C): miscible
- Melting point: -99 °C
- Boiling point: 65 °C
- Flash pt. 11 °C
- Ignition temp.: 455 °C
- LD 50 (oral, rat): 5628 mg/kg
- ADR: 3 FT1 II UN 1230
- IMDG: 3 II UN 1230
- IATA/ICAO: 3 II UN 1230
- GHS-signal word: Danger
- GHS-H sentences: H225 - H301 - H311 - H331 - H370
- GHS-P sentences: P210 - P241 - P303+P361+P353 - P370+P378 - P405 - P501a
- Tariff number: 2845 90 10 00
- Applications: for nuclear magnetic resonance spectroscopy.

SPECIFICATIONS

deuteration degree: min. 99,8 %

water (K.F., H₂O + D₂O): max. 0,03 %

performance test (NMR-spectrum): passes test



Pyridine-d₅, deuteration degree min. 99,95%, NMR spectroscopy grade, Spectrosol®



Tetrahydrofuran-d₈, deuteration degree min. 99,5%, NMR spectroscopy grade, Spectrosol®



Toluene-d₈, deuteration degree min. 99,5%, NMR spectroscopy grade, Spectrosol®



Trifluoroacetic acid-d, deuteration degree min. 99,5%, NMR spectroscopy grade, Spectrosol®

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