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эл.почта: tuf@nt-rt.ru || сайт: https://tianfu.nt-rt.ru

1,4-Benzoquinone dioxime CAS 105-11-3

Name: 1,4-Benzoquinone dioxime CAS No.:105-11-3 Molecular Structure: Molecular Structure of 105-11-3 (1,4-Benzoquinone dioxime) Formula: C6H6N2O2 Molecular Weight : 138.12 Synonyms: 2,5-Cyclohexadiene-1,4-dione,dioxime (9CI);Benzoquinone dioxime (6CI);p-Benzoquinone, dioxime (8CI);Actor Q;GM-P;NSC 14433;NSC

Mometasone CAS 105102-22-5

Name: Pregna-1,4-diene-3,20-dione,9,21-dichloro-11,17-dihydroxy-16-methyl-,(57264443,11b,16a)- CAS No.:105102-22-5 Molecular Structure: Molecular Structure of 105102-22-5 (Pregna-1,4-diene-3,20-dione,9,21-dichloro-11,17-dihydroxy-16-methyl-,(57264444,11b,16a)-) Formula: C22H28Cl2O4 Molecular Weight : 521.44

O-2-Naphthyl chlorothioformate CAS 10506-37-3

Name: Carbonochloridothioicacid, O-2-naphthalenyl ester CAS No.:10506-37-3 Molecular Structure: Molecular Structure of 10506-37-3 (Carbonochloridothioicacid, O-2-naphthalenyl ester) Formula: C11H7ClOS Molecular Weight : 222.69 Synonyms: Formicacid, chlorothio-, O-2-naphthyl ester (7CI,8CI);O-(b-Naphthyl) thiocarbonic

4-METHOXYMANDELIC ACID CAS 10502-44-0

Name: Benzeneacetic acid, alpha-hydroxy-4-methoxy- CAS No.:10502-44-0 Molecular Structure: Molecular Structure of 10502-44-0 (Benzeneacetic acid, alpha-hydroxy-4-methoxy-) Formula: C9H10O4 Molecular Weight : 182.17 Synonyms: Mandelicacid, p-methoxy- (6CI,7CI,8CI);2-Hydroxy-2-(4-methoxyphenyl)acetic

2,3-Cyclohexeno pyridine CAS 10500-57-9

Name: 2,3-Cyclohexenopyridine CAS No.:10500-57-9 Molecular Structure: Molecular Structure of 10500-57-9 (2,3-Cyclohexenopyridine) Formula: C9H11N Molecular Weight : 133.19 Synonyms: 5.6.7.8-tetra-hydroquinoline;Quinoline, 5,6,7,8-tetrahydro-;2,3-Cyclohexenopyridine;5,6,7,8-Tetra-hydroquinoline;2,3-Cyclohexeno

-Benzoyl-5'-O-(4,4-Dimethoxytrityl)-2'-O-[(tert-butyl)dimethylsilyl]adenosine-3'-(2-cyanoethyl-N,N-diisopropyl)phosphoramidite CAS 104992-55-4

Name: Adenosine,N-benzoyl-5'-O-[bis(4-methoxyphenyl)phenylmethyl]-2'-O-[(1,1-dimethylethyl)dimethylsilyl]-,3'-[2-cyanoethyl bis(1-methylethyl)phosphoramidite] (9CI) CAS No.:104992-55-4 Molecular Structure: Molecular Structure of 104992-55-4

LithiuM 2-(2', 2''-bipyridine-6'-yl)phenolate CAS 1049805-81-3

Name: LithiuM 2-(2', 2''-bipyridine-6'-yl)phenolate CAS No.:1049805-81-3 Molecular Structure: Molecular Structure of 1049805-81-3 (LithiuM 2-(2', 2''-bipyridine-6'-yl)phenolate) Formula: Molecular Weight : 0 Synonyms: LithiuM 2-(2', 2''-bipyridine-6'-yl)phenolate;Libpp EINECS: -0 Density: Melting Point: Boiling Point:

4-(Methylmercapto)aniline CAS 104-96-1

Name: Benzenamine,4-(methylthio)- CAS No.:104-96-1 Molecular Structure: Molecular Structure of 104-96-1 (Benzenamine,4-(methylthio)-) Formula: C7H9NS Molecular Weight : 139.23 Synonyms: Aniline,p-(methylthio)-

6-BROMO-2,2'-BIPYRIDINE CAS 10495-73-5

Name: 6-Bromo-2,2'-bipyridine CAS No.:10495-73-5 Molecular Structure: Molecular Structure of 10495-73-5 (6-Bromo-2,2'-bipyridine) Formula: C10H7BrN2 Molecular Weight : 235.08 Synonyms: 2-Bromo-6-(2-pyridyl)pyridine;2,2'-Bipyridine,6-bromo-;6-Bromo-2,2'-bipyridyl;6-Bromo-[2,2']bipyridinyl; EINECS: Density: 1.494 g/cm3

1H-Purin-2-amine, 6-azido- CAS 10494-88-9

Name: 1H-Purin-2-amine, 6-azido- CAS No.:10494-88-9 Molecular Structure: Molecular Structure of 10494-88-9 (1H-Purin-2-amine, 6-azido-) Formula: C5H4N8 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information:

4-Bromoanisole CAS 104-92-7

Name: 4-Bromoanisole CAS No.:104-92-7 Molecular Structure: Molecular Structure of 104-92-7 (4-Bromoanisole) Formula: C7H7BrO Molecular Weight : 187.04 Synonyms: 4-Methoxyphenyl bromide;1-bromo-4-methoxy-benzene;1-Bromo-4-methoxybenzene;p-Bromanisole;p-Methoxyphenyl bromide;Anisyl

4-Nitrosophenol CAS 104-91-6

Name: Phenol, 4-nitroso- CAS No.:104-91-6 Molecular Structure: Molecular Structure of 104-91-6 (Phenol, 4-nitroso-) Formula: C6H5NO2 Molecular Weight : 123.11 Synonyms: Phenol,p-nitroso- (6CI,8CI);4-Nitrosophenol;NSC 3124;Nitrosophenol;p-Nitrosophenol; EINECS: 203-251-6 Density: 1.236 g/cm3 Melting Point: Boiling

Rhodium(III) sulfate CAS 10489-46-0

Name: Sulfuric acid,rhodium(3+) salt (3:2) CAS No.:10489-46-0 Molecular Structure: Molecular Structure of 10489-46-0 (Sulfuric acid,rhodium(3+) salt (3:2)) Formula: Rh2S3O12 Molecular Weight : 494.00 Synonyms: Rhodiumsulfate (6CI,7CI);Dirhodium trisulfate;Rhodex;Rhodium sulfate (Rh2(SO4)3); EINECS: 234-014-5 Density:

Sodium perborate tetrahydrate CAS 10486-00-7

Name: Perboric acid(HBO(O2)), sodium salt, hydrate (1:1:4) CAS No.:10486-00-7 Molecular Structure: Molecular Structure of 10486-00-7 (Perboric acid(HBO(O2)), sodium salt, hydrate (1:1:4)) Formula: BHO3.4H2O.Na Molecular Weight : 154.86 Synonyms: Perboricacid (HBO(O2)), sodium salt, tetrahydrate (9CI);Perboric acid

p-Tolunitrile CAS 104-85-8

Name: Benzonitrile,4-methyl- CAS No.:104-85-8 Molecular Structure: Molecular Structure of 104-85-8 (Benzonitrile,4-methyl-) Formula: C8H7N Molecular Weight : 117.15 Synonyms: 1-Cyano-4-methylbenzene;4-Cyanotoluene;4-Methylbenzonitrile;4-Methylcyanobenzene;4-Methylphenyl cyanide;4-Toluenitrile;4-Tolyl cyanide;NSC

p-Tolunitrile CAS 10485-09-3

Name: 1H-Indene, 2-bromo- CAS No.:10485-09-3 Molecular Structure: Molecular Structure of 10485-09-3 (1H-Indene, 2-bromo-) Formula: C9H7Br Molecular Weight : 195.06 Synonyms: Indene,2-bromo- (6CI,8CI);2-Bromo-1H-indene;2-Indenyl bromide; EINECS: Density: 1.57 g/cm3 Melting Point: Boiling Point: 256.9 °C at 760 mmHg

4-Methylbenzylamine CAS 104-84-7

Name: Benzenemethanamine,4-methyl- CAS No.:104-84-7 Molecular Structure: Molecular Structure of 104-84-7 (Benzenemethanamine,4-methyl-) Formula: C8H11N Molecular Weight : 121.18 Synonyms: Benzylamine,p-methyl- (8CI);((4-Methylphenyl)methyl)amine;4-Aminomethyltoluene;4-Methylbenzenemethanamine;NSC

ALLYL SALICYLATE CAS 10484-09-0

Name: Benzoic acid,2-hydroxy-, 2-propen-1-yl ester CAS No.:10484-09-0 Molecular Structure: Molecular Structure of 10484-09-0 (Benzoic acid,2-hydroxy-, 2-propen-1-yl ester) Formula: C10H10 O3 Molecular Weight : 178.18 Synonyms: Benzoicacid, 2-hydroxy-, 2-propenyl ester (9CI); Salicylic acid, allyl ester(6CI,8CI); Allyl

2'-HYDROXYACETOPHENONE CAS 104809-67-8

Name: Ethanone,1-(2-hydroxyphenyl)-, labeled with carbon-14 (9CI) CAS No.:104809-67-8 Molecular Structure: Molecular Structure of 104809-67-8 (Ethanone,1-(2-hydroxyphenyl)-, labeled with carbon-14 (9CI)) Formula: C8H8 O2 Molecular Weight : 136.15 Synonyms: FEMA 3548;AKOS 90575;AKOS BBS-00003239;2-HYDROXYPHENYL METHYL

ALPHA-D(+)MALTOSE 1-PHOSPHATE DIPOTASSIUM SALT CAS 104808-98-2

Name: a-D-Glucopyranose, 4-O-a-D-glucopyranosyl-, 1-(dihydrogenphosphate), dipotassium salt (9CI) CAS No.:104808-98-2 Molecular Structure: Molecular Structure of 104808-98-2 (a-D-Glucopyranose, 4-O-a-D-glucopyranosyl-, 1-(dihydrogenphosphate), dipotassium salt (9CI)) Formula: C12H23 O14 P . 2 K Molecular Weight :

AB 1010 Mesylate CAS 1048007-93-7

Name: Masitinib (methanesulfonate) CAS No.:1048007-93-7 Molecular Structure: Molecular Structure of 1048007-93-7 (Masitinib (methanesulfonate)) Formula: C29H34N6O4S2 Molecular Weight : 594.74800 Synonyms: 4-[(4-methylpiperazin-1-yl)methyl]-N-(4-methyl-3-{[4-(pyridin-3-yl)-1,3-thiazol-2-yl]amino}phenyl)benzamide

99% CAS 104778-05-4

Name: Kazinol J CAS No.:104778-05-4 Molecular Structure: Molecular Structure of 104778-05-4 (Kazinol J) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

GSK-2110183C CAS 1047644-62-1

Name: GSK-2110183C CAS No.:1047644-62-1 Molecular Structure: Molecular Structure of 1047644-62-1 (GSK-2110183C) Formula: Molecular Weight : 0 Synonyms: Afuresertib;GSK

Dodecylpyridinium chloride CAS 104-74-5

Name: Pyridinium, 1-dodecyl-,chloride (1:1) CAS No.:104-74-5 Molecular Structure: Molecular Structure of 104-74-5 (Pyridinium, 1-dodecyl-,chloride (1:1)) Formula: C17H30N.Cl Molecular Weight : 283.93 Synonyms: 1-Dodecylpyridiniumchloride (6CI);Pyridinium, 1-dodecyl-, chloride

2,2-Bis(hydroxymethyl)propionic acid CAS 1047217

0.99

Quinacridone CAS 1047-16-1

Name: Pigment Violet 19 CAS No.:1047-16-1 Molecular Structure: Molecular Structure of 1047-16-1 (Pigment Violet 19) Formula: C20H12N2O2 Molecular Weight : 312.32 Synonyms: Cinquasia Violet R;Dark violet;Fastogen Red 7094Y;Hostaperm Red E 2B70;Hostaperm Red E 3B;HostapermRed E 5B;Hostaperm Red E 5B02;Hostaperm Red E

GERANYLGERANYL DIPHOSPHATE, TRISAMMONIUM SALT CAS 104715-21-1

Name: Diphosphoric acid, mono(3,7,11,15-tetramethylhexadecyl) ester, ion(3-) (9CI) CAS No.:104715-21-1 Molecular Structure: Molecular Structure of 104715-21-1 (Diphosphoric acid, mono(3,7,11,15-tetramethylhexadecyl) ester, ion(3-) (9CI)) Formula: C20H41O7P2 Molecular Weight : 501.534722 Synonyms: GERANYLGERANYL

Potassium perrhenate CAS 10466-65-6

Name: Rhenate (ReO41-),potassium (1:1),(57264412,T-4)- CAS No.:10466-65-6 Molecular Structure: Molecular Structure of 10466-65-6 (Rhenate (ReO41-),potassium (1:1),(57264413,T-4)-) Formula: KReO4 Molecular Weight : 289.30 Synonyms: Perrhenicacid (HReO4), potassium salt (8CI);Potassium perrhenate

L-Leucinamide hydrochloride CAS 10466-61-2

Name: Pentanamide,2-amino-4-methyl-, hydrochloride (1:1),(57264407,2S)- CAS No.:10466-61-2 Molecular Structure: Molecular Structure of 10466-61-2 (Pentanamide,2-amino-4-methyl-, hydrochloride (1:1),(57264408,2S)-) Formula: C6H14N2O.HCl Molecular Weight : 166.65 Synonyms: Leucinamide,monohydrochloride, L-

99% CAS 104662-20-6

Name: 9-Anthracenecarboxaldehyde, 2-hydroxy- CAS No.:104662-20-6 Molecular Structure: Molecular Structure of 104662-20-6 (9-Anthracenecarboxaldehyde, 2-hydroxy-) Formula: C15H10O2 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk

Pramipexole CAS 104632-26-0

Name: Pramipexole CAS No.:104632-26-0 Molecular Structure: Molecular Structure of 104632-26-0 (Pramipexole) Formula: C10H17N3S Molecular Weight : 211.33 Synonyms: Pramipexole(S)-2-Amino-6-propylamino-4,5,6,7-tetrahydrobenzthiazole; EINECS: Density: 1.17 g/cm3 Melting Point: Boiling Point: 378 °C at 760 mmHg Flash

N-(4-Oxocyclohexyl)phthalimide CAS 104618-32-8

Name: 1H-Isoindole-1,3(2H)-dione,2-(4-oxocyclohexyl)- CAS No.:104618-32-8 Molecular Structure: Molecular Structure of 104618-32-8 (1H-Isoindole-1,3(2H)-dione,2-(4-oxocyclohexyl)-) Formula: C14H13NO3 Molecular Weight : 243.26 Synonyms:

Anaritide Acetate CAS 104595-79-1

Name: Anaritide acetate [USAN] CAS No.:104595-79-1 Molecular Structure: Molecular Structure of 104595-79-1 (Anaritide acetate [USAN]) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

Phenethyl caffeate CAS 104594-70-9

Name: Phenethyl caffeate CAS No.:104594-70-9 Molecular Structure: Molecular Structure of 104594-70-9 (Phenethyl caffeate) Formula: C17H16O4 Molecular Weight : 284.31 Synonyms: 2-Phenylethylcaffeate;2-Propenoic acid,3-(3,4-dihydroxyphenyl)-, 2-phenylethyl ester;2-Phenylethyl 3-(3,4-dihydroxyphenyl)-2-propenoate;

Menthone CAS 10458-14-7

Name: Cyclohexanone,5-methyl-2-(1-methylethyl)- CAS No.:10458-14-7 Molecular Structure: Molecular Structure of 10458-14-7 (Cyclohexanone,5-methyl-2-(1-methylethyl)-) Formula: C10H18O Molecular Weight : 154.25 Synonyms:

2-Deoxy-L-ribose-anilide CAS 104578-89-4

Name: L-erythro-Pentofuranosylamine,2-deoxy-N-phenyl- CAS No.:104578-89-4 Molecular Structure: Molecular Structure of 104578-89-4 (L-erythro-Pentofuranosylamine,2-deoxy-N-phenyl-) Formula: C11H15 N O3 Molecular Weight : 0 Synonyms: 2-Deoxy-L-ribose-anilide EINECS: Density: 1.305 Melting Point: Boiling Point: Flash

Benzyl formate CAS 104-57-4

Name: Formic acid,phenylmethyl ester CAS No.:104-57-4 Molecular Structure: Molecular Structure of 104-57-4 (Formic acid,phenylmethyl ester) Formula: C8H8O2 Molecular Weight : 136.15 Synonyms: Formicacid, benzyl ester (6CI,7CI,8CI);Benzyl alcohol, formate;Benzyl formate;Benzyl methanoate;NSC 8049; EINECS: 203-214-4

1-Chloro-3-phenylpropane CAS 104-52-9

Name: Benzene,(57264396,3-chloropropyl)- CAS No.:104-52-9 Molecular Structure: Molecular Structure of 104-52-9 (Benzene,(57264397,3-chloropropyl)-) Formula: C9H11Cl Molecular Weight : 154.64 Synonyms: g-Chloropropylbenzene;g-Phenylpropyl chloride;NSC

(3-QUINUCLIDINYL)DIPHENYL CARBINOL HYDROCHLORIDE CAS 10447-39-9

Name: 1-Azabicyclo[2.2.2]octane-3-methanol,a,a-diphenyl- CAS No.:10447-39-9 Molecular Structure: Molecular Structure of 10447-39-9 (1-Azabicyclo[2.2.2]octane-3-methanol,a,a-diphenyl-) Formula: C20H23NO Molecular Weight : 329.86 Synonyms: 3-Quinuclidinemethanol,a,a-diphenyl- (7CI,8CI);Diphenyl-3-quinuclidylcarbinol;

4-Methoxybenzyl cyanide CAS 104-47-2

Name: Benzeneacetonitrile,4-methoxy- CAS No.:104-47-2 Molecular Structure: Molecular Structure of 104-47-2 (Benzeneacetonitrile,4-methoxy-) Formula: C9H9NO Molecular Weight : 147.17 Synonyms: Acetonitrile,(57264393,p-methoxyphenyl)-

CERAMIDE CAS 104404-17-3

Name: Ceramide CAS No.:104404-17-3 Molecular Structure: Molecular Structure of 104404-17-3 (Ceramide) Formula: Molecular Weight : 536.89 Synonyms: CERAMIDE;Ceramide?(saturated,?unsaturated?and?a-OH?fatty?acids);Ceramide (d18:1/12:0);N-((2S,3R,E)-1,3-dihydroxyoctadec-4-en-2-yl)dodecanamide;Ceramide

Poly(oxy-1,2-ethanediyl), .alpha.-1,1-biphenyl-4-yl-.omega.-hydroxy-, benzylated CAS 104376-72-9

Name: Poly(oxy-1,2-ethanediyl), α-[1,1'-biphenyl]-4-yl-ω-hydroxy-, benzylated CAS No.:104376-72-9 Molecular Structure: Molecular Structure of 104376-72-9 (Poly(oxy-1,2-ethanediyl), α-[1,1'-biphenyl]-4-yl-ω-hydroxy-, benzylated) Formula: Unspecified Molecular Weight : 0 Synonyms: Poly(oxy-1,2-ethanediyl),

2-Propenoic acid, 3-[(4'-cyano[1,1'-biphenyl]-4-yl)oxy]propyl ester CAS 104357-57-5

Name: 2-Propenoic acid, 3-[(4'-cyano[1,1'-biphenyl]-4-yl)oxy]propyl ester CAS No.:104357-57-5 Molecular Structure: Molecular Structure of 104357-57-5 (2-Propenoic acid, 3-[(4'-cyano[1,1'-biphenyl]-4-yl)oxy]propyl ester) Formula: Molecular Weight : 307.34318 Synonyms: 2-Propenoic acid,

O-BENZYL-D-SERINE CAS 10433-52-0

Name: D-Serine,O-(phenylmethyl)- CAS No.:10433-52-0 Molecular Structure: Molecular Structure of 10433-52-0 (D-Serine,O-(phenylmethyl)-) Formula: C10H13NO3 Molecular Weight : 195.22 Synonyms: Alanine,3-(benzyloxy)-, D- (8CI);O-Benzyl-D-serine; EINECS: 233-916-6 Density: 1.218 g/cm3 Melting Point: Boiling Point: 359.053

: 2-Ethyl-2-oxazoline CAS 10431-98-8

Name: Oxazole,2-ethyl-4,5-dihydro- CAS No.:10431-98-8 Molecular Structure: Molecular Structure of 10431-98-8 (Oxazole,2-ethyl-4,5-dihydro-) Formula: C5H9NO Molecular Weight : 99.13 Synonyms: 2-Oxazoline,2-ethyl-(6CI,7CI,8CI);2-Ethyl-2-oxazoline;2-Ethyl-4,5-dihydrooxazole;2-Ethyloxazoline;NSC 136557; EINECS: 233-912-4

Benzonatate CAS 104-31-4

Name: Benzoic acid,4-(butylamino)-, 3,6,9,12,15,18,21,24,27-nonaoxaoctacos-1-yl ester CAS No.:104-31-4 Molecular Structure: Molecular Structure of 104-31-4 (Benzoic acid,4-(butylamino)-, 3,6,9,12,15,18,21,24,27-nonaoxaoctacos-1-yl ester) Formula: C30H53NO11 Molecular Weight : 603.75 Synonyms: Benzoicacid,

dithieno[3,2-c:2',3'-e]oxepine-4,6-dione CAS 1043023-52-4

Name: dithieno[3,2-c:2',3'-e]oxepine-4,6-dione CAS No.:1043023-52-4 Molecular Structure: Molecular Structure of 1043023-52-4 (dithieno[3,2-c:2',3'-e]oxepine-4,6-dione) Formula: Molecular Weight : 236.26696 Synonyms: dithieno[3,2-c:2',3'-e]oxepine-4,6-dione EINECS: Density: Melting Point: Boiling Point: Flash Point:

(1R,2R,4S)-2-[(5-hexen-1-ylmethylamino)carbonyl]-4-[[7-methoxy-8-methyl- 2-[4-(1-isopropyl)-2-thiazolyl]-4-quinolinyl]oxy]-Cyclopentanecarboxylic acid methyl ester CAS 1042695-87-3

Name: (1R,2R,4S)-2-[(5-hexen-1-ylmethylamino)carbonyl]-4-[[7-methoxy-8- methyl- 2-[4-(1-isopropyl)-2-thiazolyl]-4-quinolinyl]oxy]- Cyclopentanecarboxylic acid methyl ester CAS No.:1042695-87-3 Molecular

Structure: Molecular Structure of 1042695-87-3 ((1R,2R,4S)-2-[(5-hexen-1-ylmethylamino)carbonyl]-4-[[7-methoxy-8-

3,4-Dihydroxybenzophenone CAS 10425-11-3

Name: 3,4-Dihydroxybenzophenone CAS No.:10425-11-3 Molecular Structure: Molecular Structure of 10425-11-3 (3,4-Dihydroxybenzophenone) Formula: C13H10O3 Molecular Weight : 214.2167 Synonyms: Methanone,(57264382,3,4-dihydroxyphenyl)phenyl-;Benzophenone,3,4-dihydroxy- (7CI,8CI); EINECS: Density: 1.302 g/cm3 Melting

1-(5-Bromo-2-methoxy-phenyl)adamantane CAS 104224-63-7

Name: 1-(5-Bromo-2-methoxy-phenyl)adamantane CAS No.:104224-63-7 Molecular Structure: Molecular Structure of 104224-63-7 (1-(5-Bromo-2-methoxy-phenyl)adamantane) Formula: C17H21BrO Molecular Weight : 321.25 Synonyms: 1-(5-Bromo-2-methoxyphenyl)-tricyclo[3.3.1.13,7]decane;2-Adamantyl-4-bromoanisole; EINECS: Density:

AMBERLITE XAD-16 CAS 104219-63-8

Name: Amberlite xad-16 CAS No.:104219-63-8 Molecular Structure: Molecular Structure of 104219-63-8 (Amberlite xad-16) Formula: Unspecified Molecular Weight : 0 Synonyms: Amberlite XAD-16 EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols:R36:; Risk Codes: Xi:; Transport

2-Chloromandelic acid CAS 10421-85-9

Name: Benzeneacetic acid,2-chloro-a-hydroxy- CAS No.:10421-85-9 Molecular Structure: Molecular Structure of 10421-85-9 (Benzeneacetic acid,2-chloro-a-hydroxy-) Formula: C8H7ClO3 Molecular Weight : 186.59 Synonyms: DL-2-Chloromandelic acid;Mandelicacid, o-chloro- (6CI,7CI,8CI);(2-Chlorophenyl)glycolic

Dimethyl acetylsuccinate CAS 10420-33-4

Name: Butanedioic acid,2-acetyl-, 1,4-dimethyl ester CAS No.:10420-33-4 Molecular Structure: Molecular Structure of 10420-33-4 (Butanedioic acid,2-acetyl-, 1,4-dimethyl ester) Formula: C8H12O5 Molecular Weight : 188.18 Synonyms: Butanedioicacid, acetyl-, dimethyl ester (9CI);Succinic acid, acetyl-, dimethyl

3,5-DIFLUORO-4-METHOXYBENZONITRILE CAS 104197-15-1

Name: 3,5-Difluoro-4-methoxybenzonitrile CAS No.:104197-15-1 Molecular Structure: Molecular Structure of 104197-15-1 (3,5-Difluoro-4-methoxybenzonitrile) Formula: C8H5F2NO Molecular Weight : 169.13 Synonyms: 3,5-Difluoro-4-methoxybenzonitrile; EINECS: Density: 1.279 g/cm3 Melting Point: Boiling Point: 250.592 °C at

TETRAPHENYLARSONIUM CHLORIDE MONOHYDRATE CAS 104170-16-3

Name: Arsonium, tetraphenyl-,chloride, hydrate (1:1:1) CAS No.:104170-16-3 Molecular Structure: Molecular Structure of 104170-16-3 (Arsonium, tetraphenyl-,chloride, hydrate (1:1:1)) Formula: C24H20 As . Cl . H2 O Molecular Weight : 436.81 Synonyms: Arsonium,tetraphenyl-, chloride, monohydrate (9CI) EINECS: Density:

Disperse Blue 354 CAS 104137-27-1

Name: Disperse Blue 354 CAS No.:104137-27-1 Molecular Structure: Molecular Structure of 104137-27-1 (Disperse Blue 354) Formula: C31H37N3O2S Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

： 6aH-Cyclohepta[a]naphthalene-6a,9-dicarboxaldehyde,1,2,3,4,4a,5,6,7,8,11,11a,11b-dodecahydro-7-hydroxy-4,4,11b-trimethyl-,(57264370,4aS,6aS,7R,11aR,11bS)- CAS 104113-52-2

Name: 6aH-Cyclohepta[a]naphthalene-6a,9-dicarboxaldehyde,1,2,3,4,4a,5,6,7,8,11,11a,11b-dodecahydro-7-hydroxy-4,4,11b-trimethyl-,(57264371,4aS,6aS,7R,11aR,11bS)- CAS No.:104113-52-2 Molecular Structure: Molecular Structure of 104113-52-2

99% CAS 10411-24-2

0.99

3,5-Diiodo-L-thyronine CAS 1041-01-6

Name: 3,5-Diiodo-L-thyronine CAS No.:1041-01-6 Molecular Structure: Molecular Structure of 1041-01-6 (3,5-Diiodo-L-thyronine) Formula: C15H13I2NO4 Molecular Weight : 525.077 Synonyms: 3,5-Diiodothyronine (VAN);L-Tyrosine, O- (4-hydroxyphenyl)-3,5-diiodo-;Diiodo-L-thyronine;L-Tyrosine, O-(4-hydroxyphenyl)-3,5-diiodo-

Fmoc-D-glutamic acid gamma-tert-butyl ester CAS 104091-08-9

Name: D-Glutamicacid, N-[(9H-fluoren-9-ylmethoxy)carbonyl]-, 5-(1,1-dimethylethyl) ester CAS No.:104091-08-9 Molecular Structure: Molecular Structure of 104091-08-9 (D-Glutamicacid, N-[(9H-fluoren-9-ylmethoxy)carbonyl]-, 5-(1,1-dimethylethyl) ester) Formula: C24H27NO6 Molecular Weight : 425.48 Synonyms:

2,3-dihydro-1H-inden-2-yl(Methyl)aMine(HCl) CAS 10408-85-2

Name: 2,3-dihydro-1H-inden-2-yl(methyl)amine(HCl) CAS No.:10408-85-2 Molecular Structure: Molecular Structure of 10408-85-2 (2,3-dihydro-1H-inden-2-yl(methyl)amine(HCl)) Formula: Molecular Weight : 183.68 Synonyms: N-Methyl-2-indanamine hydrochloride

METHYL 3-METHOXY-2-PENTENOATE CAS 104065-67-0

Name: 2-Pentenoic acid,3-methoxy-, methyl ester CAS No.:104065-67-0 Molecular Structure: Molecular Structure of 104065-67-0 (2-Pentenoic acid,3-methoxy-, methyl ester) Formula: C7H12 O3 Molecular Weight : 144.17 Synonyms: Methyl3-methoxy-2-pentenoate; Methyl 3-methoxy-3-ethylacrylate EINECS: Density: 0.978g/cm3

Trospium chloride CAS 10405-02-4

Name: Trospium chloride CAS No.:10405-02-4 Molecular Structure: Molecular Structure of 10405-02-4 (Trospium chloride) Formula: C25H30NO3 Molecular Weight : 427.96 Synonyms: Spiro[1aH,5aH-nortropane-8,1'-pyrrolidinium], 3a-hydroxy-, chloride,

99% CAS 1040375-91-4

0.99

2-Bromo-5-nitroaniline CAS 10403-47-1

Name: Benzenamine,2-bromo-5-nitro- CAS No.:10403-47-1 Molecular Structure: Molecular Structure of 10403-47-1 (Benzenamine,2-bromo-5-nitro-) Formula: C6H5BrN2O2 Molecular Weight : 217.02 Synonyms: Aniline,2-bromo-5-nitro- (6CI,7CI,8CI);(2-Bromo-5-nitrophenyl)amine;2-Bromo-5-nitrobenzenamine; EINECS: 233-874-9 Density:

4-Nitrophenylacetic acid CAS 104-03-0

Name: Benzeneaceticacid, 4-nitro- CAS No.:104-03-0 Molecular Structure: Molecular Structure of 104-03-0 (Benzeneaceticacid, 4-nitro-) Formula: C8H7NO4 Molecular Weight : 181.15 Synonyms: Aceticacid,(57264359,p-nitrophenyl)- (6CI,8CI);(4-Nitrophenyl)acetic acid;(p-Nitrophenyl)acetic acid;2-(4-Nitrophenyl)acetic

OlMesartan DiMer Ester IMpurity CAS 1040250-19-8

Name: Olmesartan Dimer Ester Impurity CAS No.:1040250-19-8 Molecular Structure: Molecular Structure of 1040250-19-8 (Olmesartan Dimer Ester Impurity) Formula: Molecular Weight : 872.97212 Synonyms: OlMesartan DiMer Ester

RICINOLEIC ACID N-HEXADECYL ESTER CAS 10401-55-5

Name: 9-Octadecenoic acid,12-hydroxy-, hexadecyl ester,(57264355,9Z,12R)- CAS No.:10401-55-5 Molecular Structure: Molecular Structure of 10401-55-5 (9-Octadecenoic acid,12-hydroxy-, hexadecyl ester,(57264356,9Z,12R)-) Formula: C34H66O3 Molecular Weight : 522.89 Synonyms: 9-Octadecenoicacid, 12-hydroxy-, hexadecyl

NICOTINYL CHLORIDE HYDROCHLORIDE CAS 10400-19-8

Name: 3-Pyridinecarbonylchloride CAS No.:10400-19-8 Molecular Structure: Molecular Structure of 10400-19-8 (3-Pyridinecarbonylchloride) Formula: C6H4ClNO Molecular Weight : 141.5551 Synonyms: 3-Pyridinecarboxylic acid chloride;3-Pyridoyl chloride;3-Pyridylcarbonyl

Phenylacetone CAS 103-97-7

0.99

p-Toluenesulfonyl semicarbazide CAS 10396-10-8

Name: p-Toluenesulfonyl semicarbazide CAS No.:10396-10-8 Molecular Structure: Molecular Structure of 10396-10-8 (p-Toluenesulfonyl semicarbazide) Formula: C8H11N3O3S Molecular Weight : 229.26 Synonyms: Benzenesulfonic acid,4-methyl-, 2-(aminocarbonyl)hydrazide;Semicarbazide,1-(p-tolylsulfonyl)-

(4alpha)-2beta,15alpha-dihydroxy-19-norkaur-16-en-18-oic acid CAS 10391-47-6

Name: 19-Norkaur-16-en-18-oicacid, 2,15-dihydroxy-,(57264349,2b,4a,15a)- CAS No.:10391-47-6 Molecular Structure: Molecular Structure of 10391-47-6 (19-Norkaur-16-en-18-oicacid, 2,15-dihydroxy-,(57264350,2b,4a,15a)-) Formula: C19H28 O4 Molecular Weight : 320.4232 Synonyms: 19-Norkaur-16-en-18-oicacid,

Maxacalcitol CAS 103909-75-7

Name: Maxacalcitol CAS No.:103909-75-7 Molecular Structure: Molecular Structure of 103909-75-7 (Maxacalcitol) Formula: C26H42O4 Molecular Weight : 418.6093 Synonyms: 1,3-Cyclohexanediol,4-methylene-5-[(2E)-[(1S,3aS,7aS)-octahydro-1-[(1S)-1-(3-hydroxy-3-methylbutoxy)ethyl]-7a-methyl-4H-inden-4-ylidene]ethylidene]-,(

TRI-P-TOLYLPHOSPHINE CAS 1038-95-5

Name: Phosphine,tris(4-methylphenyl)- CAS No.:1038-95-5 Molecular Structure: Molecular Structure of 1038-95-5 (Phosphine,tris(4-methylphenyl)-) Formula: C21H21P Molecular Weight : 304.37 Synonyms: Phosphine,tri-p-tolyl- (6CI,7CI,8CI);NSC 97371;TPTP;TPTP

4'-Bromoacetanilide CAS 103-88-8

Name: Acetamide,N-(4-bromophenyl)- CAS No.:103-88-8 Molecular Structure: Molecular Structure of 103-88-8 (Acetamide,N-(4-bromophenyl)-) Formula: C8H8BrNO Molecular Weight : 214.08 Synonyms: Acetanilide,4'-bromo- (6CI,7CI,8CI);Acetanilide, p-bromo-

Nsc267697 CAS 10387-50-5

Name: 2H-1-Benzopyran-2-one,7-[(3-methyl-2-buten-1-yl)oxy]- CAS No.:10387-50-5 Molecular Structure: Molecular Structure of 10387-50-5 (2H-1-Benzopyran-2-one,7-[(3-methyl-2-buten-1-yl)oxy]-) Formula: C14H14 O3 Molecular Weight : 230.2592 Synonyms: 2H-1-Benzopyran-2-one,7-[(3-methyl-2-butenyl)oxy]- (9CI); Coumarin,

1-PHENYL-2-THIOUREA CAS 103-85-5

Name: 1-Phenyl-2-thiourea CAS No.:103-85-5 Molecular Structure: Molecular Structure of 103-85-5 (1-Phenyl-2-thiourea) Formula: C7H8N2S Molecular Weight : 152.23 Synonyms: Thiourea,phenyl- (9CI);Urea, 1-phenyl-2-thio- (8CI);1-Phenyl-2-thiourea;1-Phenylthiourea;N-Phenylthiourea;NSC

99% CAS 10-38-4

0.99

NICKEL (II) PHOSPHATE CAS 10381-36-9

Name: Phosphoric acid,nickel(2+) salt (2:3) CAS No.:10381-36-9 Molecular Structure: Molecular Structure of 10381-36-9 (Phosphoric acid,nickel(2+) salt (2:3)) Formula: H3O4 P . 3/2 Ni Molecular Weight : 366.02 Synonyms: Nickelphosphate (Ni3(PO4)2) (6CI,7CI); Nickel phosphate; Nickel phosphate (Ni3P2O8);Nickel(2+)

2-Phenylacetamide CAS 103-81-1

Name: 2-Phenylacetamide CAS No.:103-81-1 Molecular Structure: Molecular Structure of 103-81-1 (2-Phenylacetamide) Formula: C8H9NO Molecular Weight : 135.16 Synonyms: Acetamide,2-phenyl-(6CI,8CI);Benzenediacetamide (7CI);Benzeneacetamide;NSC 1877;Phenacetamide;Phenyl-b-acetylamine;Phenylacetic acid

Ethylenediaminetetraacetic acid tetrasodium salt dihydrate CAS 10378-23-1

Name: Glycine,N,N'-1,2-ethanediylbis[N-(carboxymethyl)-, sodium salt, hydrate (1:4:2) CAS No.:10378-23-1 Molecular Structure: Molecular Structure of 10378-23-1 (Glycine,N,N'-1,2-ethanediylbis[N-(carboxymethyl)-, sodium salt, hydrate (1:4:2)) Formula: C10H16N2Na4O8.2(H2O) Molecular Weight : 420.20 Synonyms:

ALLOSAMIDIN CAS 103782-08-7

Name: b-D-Allopyranoside,(57264334,3aR,4R,5R,6S,6aS)-2-(dimethylamino)-3a,5,6,6a-tetrahydro-4-hydroxy-6-(hydroxymethyl)-4H-cyclopentoxazol-5-yl2-(acetylamino)-4-O-[2-(acetylamino)-2-deoxy-b-D-allopyranosyl]-2-deoxy- CAS No.:103782-08-7 Molecular Structure: Molecular Structure of 103782-08-7 (b-D-Allopyranoside,(

LITHIUM IODIDE CAS 10377-51-2

Name: Lithium iodide CAS No.:10377-51-2 Molecular Structure: Molecular Structure of 10377-51-2 (Lithium iodide) Formula: ILi Molecular Weight : 133.85 Synonyms: Lithiumiodide (8CI);Lithium monoiodide;Lithium iodide (LiI); EINECS: 233-822-5 Density: 3.49 g/mL at 25 °C(lit.) Melting Point: Boiling Point: 1171 °C Flash

1-(2-Hydroxyethyl)piperazine CAS 103-76-4

Name: 1-(2-Hydroxyethyl)piperazine CAS No.:103-76-4 Molecular Structure: Molecular Structure of 103-76-4 (1-(2-Hydroxyethyl)piperazine) Formula: C6H14N2O Molecular Weight : 130.19 Synonyms:

2,5-Furandicarbonyldichloride(9CI) CAS 10375-34-5

Name: 2,5-Furandicarbonyldichloride CAS No.:10375-34-5 Molecular Structure: Molecular Structure of 10375-34-5 (2,5-Furandicarbonyldichloride) Formula: C6H2 Cl2 O3 Molecular Weight : 0 Synonyms: 2,5-Furandicarbonylchloride (6CI,7CI,8CI); 2,5-Furandicarboxylic acid chloride;Furan-2,5-dicarboxylic acid dichloride EINECS:

(S)-6,7-Dimethoxy-1,2,3,4-tetrahydro-3-isoquinolinecarboxylic acid hydrochloride CAS 103733-66-0

Name: 3-Isoquinolinecarboxylicacid, 1,2,3,4-tetrahydro-6,7-dimethoxy-,(57264325,3S)- CAS No.:103733-66-0 Molecular Structure: Molecular Structure of 103733-66-0 (3-Isoquinolinecarboxylicacid, 1,2,3,4-tetrahydro-6,7-dimethoxy-,(57264326,3S)-) Formula: C12H15NO4 Molecular Weight : 273.25 Synonyms:

PD 113413 CAS 103733-50-2

Name: 2H-Pyrazino[1,2-b]isoquinoline-2-aceticacid, 1,3,4,6,11,11a-hexahydro-3-methyl-1,4-dioxo-a-(2-phenylethyl)-,(57264322,aS,3S,11aS)- CAS No.:103733-50-2 Molecular Structure: Molecular Structure of 103733-50-2 (2H-Pyrazino[1,2-b]isoquinoline-2-aceticacid,

1,2,3,4-Tetrahydro-6,7-dimethoxy-3-isoquinolinecarboxylic acid phenylmethyl ester hydrochloride CAS 103733-32-0

Name: 1,2,3,4-Tetrahydro-6,7-dimethoxy-3-isoquinolinecarboxylic acid phenylmethyl ester hydrochloride CAS No.:103733-32-0 Molecular Structure: Molecular Structure of 103733-32-0 (1,2,3,4-Tetrahydro-6,7-dimethoxy-3-isoquinolinecarboxylic acid phenylmethyl ester hydrochloride) Formula: C19H21NO4.HCl Molecular Weight :

IPI-926 CAS 1037210-93-7

Name: Methanesulfonamide, N-[(2S,3R,3'R,3aS,4'aR,6S,6'aR,6'bS,7aR,12'aS,12'bS)-2',3',3a,4,4',4'a,5,5',6,6',6'a,6'b,7,7',7a,8',10',12',12'a,12'b-eicosahydro-3,6,11',12'b-tetramethylspiro[furo[3,2-b]pyridine-2(3H),9'(1'H)-naphth[2,1-a]azulen]-3'-yl]- CAS No.:1037210-93-7 Molecular Structure: Molecular Structure of

FORMANILIDE CAS 103-70-8

Name: Formamide, N-phenyl- CAS No.:103-70-8 Molecular Structure: Molecular Structure of 103-70-8 (Formamide, N-phenyl-) Formula: C7H7 N O Molecular Weight : 121.15 Synonyms: Formanilide(6CI,7CI,8CI); Carbanilaldehyde; Formamidobenzene; Formylaniline;N-Formylaniline; N-Phenylformamide; NSC 203239; NSC 8862 EINECS:

N-Methylbenzylamine CAS 103-67-3

Name: Benzenemethanamine,N-methyl- CAS No.:103-67-3 Molecular Structure: Molecular Structure of 103-67-3 (Benzenemethanamine,N-methyl-) Formula: C8H11N Molecular Weight : 121.18 Synonyms: Benzylamine,N-methyl-

1-[[4-[(4,5,6,7-Tetrachloro-3-oxo-isoindoline-1-ylidene)amino]phenyl]azo]-2-hydro xy-N-(4-methoxy-2-methylphenyl)-11H-benzo[a]carbazole-3-carboxamide CAS 103621-96-1

Name: 11H-Benzo[a]carbazole-3-carboxamide,2-hydroxy-N-(4-methoxy-2-methylphenyl)-1-[2-[4-[(4,5,6,7-tetrachloro-1-oxo-1H-isoindol-3-yl)amino]phenyl]diazenyl]- CAS No.:103621-96-1 Molecular Structure: Molecular Structure of 103621-96-1

SCANDIUM CHLORIDE CAS 10361-84-9

Name: Scandium chloride CAS No.:10361-84-9 Molecular Structure: Molecular Structure of 10361-84-9 (Scandium chloride) Formula: Cl3Sc Molecular Weight : 151.31 Synonyms: NSC 132051;Scandium chloride (ScCl3);Scandium trichloride; EINECS: 233-799-1 Density: 2.39 g/mL at 25 °C(lit.) Melting Point: Boiling Point: Flash

FEMA 2296 CAS 103-61-7

Name: Butanoic acid,3-phenyl-2-propen-1-yl ester CAS No.:103-61-7 Molecular Structure: Molecular Structure of 103-61-7 (Butanoic acid,3-phenyl-2-propen-1-yl ester) Formula: C13H16O2 Molecular Weight : 204.26 Synonyms: Butanoicacid, 3-phenyl-2-propenyl ester (9CI);Butyric acid, cinnamyl ester (6CI,8CI);Cinnamyl

BISMUTH HYDROXIDE CAS 10361-43-0

Name: Bismuth hydroxide CAS No.:10361-43-0 Molecular Structure: Molecular Structure of 10361-43-0 (Bismuth hydroxide) Formula: BiH3O3 Molecular Weight : 260.00 Synonyms: Bismon;Bismuth acid;Bismuth trihydroxide; EINECS: 233-790-2 Density: 8.9 g/cm3 Melting Point: Boiling Point: 100oC at 760 mmHg Flash Point:

AMMONIUM CARBONATE CAS 10361-29-2

Name: Carbonic acid, ammoniumsalt (1:?) CAS No.:10361-29-2 Molecular Structure: Molecular Structure of 10361-29-2 (Carbonic acid, ammoniumsalt (1:?)) Formula: CH2O3 . x H3N Molecular Weight : 79.05530 Synonyms: Carbonicacid, ammonium salt (8CI,9CI);E 503; EINECS: 233-786-0 Density: 0.780-0.830 g/m3 Melting Point:

Octadecyl gallate CAS 10361-12-3

Name: Benzoic acid,3,4,5-trihydroxy-, octadecyl ester CAS No.:10361-12-3 Molecular Structure: Molecular Structure of 10361-12-3 (Benzoic acid,3,4,5-trihydroxy-, octadecyl ester) Formula: C25H42O5 Molecular Weight : 422.60 Synonyms: Octadecylgallate;Gallicacid, octadecyl ester (6CI,7CI,8CI);1-Octadecanol, gallate

Sodium metaphosphate CAS 10361-03-2

Name: Metaphosphoric acid(HPO3), sodium salt (1:1) CAS No.:10361-03-2 Molecular Structure: Molecular Structure of 10361-03-2 (Metaphosphoric acid(HPO3), sodium salt (1:1)) Formula: NaO3P Molecular Weight : 101.96 Synonyms: METAPHOSPHORIC ACID , SODIUM SALT;METAPHOSPHORIC ACID(HPO3), SODIUM SALT (1:1);sporix;SODIUM

P005672 hydrochloride CAS 1035979-44-2

Name: (4S,4aS,5aR,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3, 10,12,12a-tetrahydroxy-7-[(methoxymethylamino)methyl]-1,11-dioxo-2-naphthacenecarboxamide hydrochloride CAS No.:1035979-44-2 Molecular Structure: Molecular Structure of 1035979-44-2

3-PHENYLPROPYL ISOBUTYRATE CAS 103-58-2

Name: Propanoic acid,2-methyl-, 3-phenylpropyl ester CAS No.:103-58-2 Molecular Structure: Molecular Structure of 103-58-2 (Propanoic acid,2-methyl-, 3-phenylpropyl ester) Formula: C13H18 O2 Molecular Weight : 206.2808 Synonyms: Isobutyric acid,3-phenylpropyl ester (7CI,8CI); 1-Propanol, 3-phenyl-, isobutyrate

Cinnamyl acetate CAS 103-54-8

Name: 2-Propen-1-ol,3-phenyl-, 1-acetate CAS No.:103-54-8 Molecular Structure: Molecular Structure of 103-54-8 (2-Propen-1-ol,3-phenyl-, 1-acetate) Formula: C11H12O2 Molecular Weight : 176.23 Synonyms: 2-Propen-1-ol,3-phenyl-, acetate (9CI);Cinnamyl alcohol, acetate

1,2-EPOXY-5-HEXENE CAS 10353-53-4

Name: 1,2-Epoxy-5-hexene CAS No.:10353-53-4 Molecular Structure: Molecular Structure of 10353-53-4 (1,2-Epoxy-5-hexene) Formula: C6H10O Molecular Weight : 98.14 Synonyms: 1-Hexene,5,6-epoxy-(7CI,8CI);Oxirane, 3-butenyl- (9CI);2-(3-Buten-1-yl)oxirane;2-(3-Butenyl)oxirane;3-Butenyloxirane;5,6-Epoxy-1-hexene;Diallyl

NA CAS 103515-22-6

0.99

Benzyl ether CAS 103-50-4

Name: Benzyl ether CAS No.:103-50-4 Molecular Structure: Molecular Structure of 103-50-4 (Benzyl ether) Formula: C14H14O Molecular Weight : 198.28 Synonyms: Benzylether (8CI);1,1'-[Oxybis(methylene)]bis[benzene];BA (plasticizer);Benzyloxide;Dibenzyl ether;NSC 5931;Plastikator BA;Benzene,1,1'-[oxybis(methylene)]bis-;

N-Cyclohexyltaurine CAS 103-47-9

Name: Ethanesulfonic acid,2-(cyclohexylamino)- CAS No.:103-47-9 Molecular Structure: Molecular Structure of 103-47-9 (Ethanesulfonic acid,2-(cyclohexylamino)-) Formula: C8H17NO3S Molecular Weight : 207.29 Synonyms: 2-[N-Cyclohexylamino]ethanesulfonic acid;CHES;N-Cyclohexyl-2-aminoethanesulfonic acid;NSC

TRIPHENYLPHOSPHINE DIBROMIDE CAS 1034-39-5

Name: Phosphorane,dibromotriphenyl- CAS No.:1034-39-5 Molecular Structure: Molecular Structure of 1034-39-5 (Phosphorane,dibromotriphenyl-) Formula: C18H15Br2P Molecular Weight : 422.09 Synonyms: NSC 87871;Triphenyldibromophosphorane;Triphenylphosphine dibromide; EINECS: 213-855-1 Density: Melting Point: Boiling

ACID YELLOW 99 CAS 10343-58-5

Name: Chromate(1-),hydroxy[2-(hydroxy-kO)-5-nitro-3-[2-[2-(oxo-kO)-1-[(phenylamino)carbonyl]propyl]diazenyl-kN1]benzenesulfonato(3-)]-, sodium (1:1),(57264298,T-4)- CAS No.:10343-58-5 Molecular Structure: Molecular Structure of 10343-58-5

2,3,4,6-Tetraacetyl-D-glucose CAS 10343-06-3(3947-62-4beta)

Name: D-Glucopyranose,2,3,4,6-tetraacetate CAS No.:10343-06-3 Molecular Structure: Molecular Structure of 10343-06-3 (D-Glucopyranose,2,3,4,6-tetraacetate) Formula: C14H20O10 Molecular Weight : 348.31 Synonyms: Glucopyranose,2,3,4,6-tetraacetate, D- (8CI);2,3,4,6-Tetra-O-acetyl-D-glucopyranose; EINECS: Density: 1.339

3',5'-di-o-p-chlorobenzoyl-2-deoxy-5-azacytosine CAS 1034301-08-0

Name: 1,3,5-Triazin-2(1H)-one,4-amino-1-[3,5-bis-O-(4-chlorobenzoyl)-2-deoxy-b-D-erythro-pentofuranosyl]- CAS No.:1034301-08-0 Molecular Structure: Molecular Structure of 1034301-08-0 (1,3,5-Triazin-2(1H)-one,4-amino-1-[3,5-bis-O-(4-chlorobenzoyl)-2-deoxy-b-D-erythro-pentofuranosyl]-) Formula: C22H18Cl2N4O6 Molecular

N,N-Diisopropylmethylamine CAS 10342-97-9

Name: 2-Propanamine,N-methyl-N-(1-methylethyl)- CAS No.:10342-97-9 Molecular Structure: Molecular Structure of 10342-97-9 (2-Propanamine,N-methyl-N-(1-methylethyl)-) Formula: C7H17 N Molecular Weight : 115.22 Synonyms: Diethylamine,N,1,1'-trimethyl- (6CI,7CI,8CI); Diisopropylmethylamine;Methyldiisopropylamine;

4'-Bromopropiophenone CAS 10342-83-3

Name: 1-Propanone,1-(4-bromophenyl)- CAS No.:10342-83-3 Molecular Structure: Molecular Structure of 10342-83-3 (1-Propanone,1-(4-bromophenyl)-) Formula: C9H9 Br O Molecular Weight : 213.0712 Synonyms: Propiophenone,4'-bromo- (6CI,7CI,8CI); 1-(4-Bromophenyl)-1-propanone;1-(4-Bromophenyl)propanone; 4-Bromophenyl ethyl

Benzyl cinnamate CAS 103-41-3

Name: 2-Propenoicacid, 3-phenyl-, phenylmethyl ester CAS No.:103-41-3 Molecular Structure: Molecular Structure of 103-41-3 (2-Propenoicacid, 3-phenyl-, phenylmethyl ester) Formula: C16H14O2 Molecular Weight : 238.29 Synonyms: Cinnamicacid, benzyl ester (6CI,7CI,8CI);3-Phenyl-2-propenoic acid benzyl ester;Benzyl

Zinc L-lactate CAS 103404-76-8

Name: Zinc L-lactate CAS No.:103404-76-8 Molecular Structure: Molecular Structure of 103404-76-8 (Zinc L-lactate) Formula: 2(C3H5O3).Zn Molecular Weight : 243.52000 Synonyms: L-Zinc Lactatel; DL-Milchsaeure,Zinklactat; zincum lactate; zinc salt of lactic acid; L-Zinc; Ammonium Salts; DL-lactic acid,zinc lactate;

Octyl gallate CAS 1034-01-1

Name: Benzoic acid,3,4,5-trihydroxy-, octyl ester CAS No.:1034-01-1 Molecular Structure: Molecular Structure of 1034-01-1 (Benzoic acid,3,4,5-trihydroxy-, octyl ester) Formula: C15H22O5 Molecular Weight : 282.33 Synonyms: Gallicacid, octyl ester (6CI,8CI);E 311;NSC 97419;Octyl 3,4,5-trihydroxybenzoate;Progallin

BENZYL ISOVALERATE CAS 103-38-8

Name: Butanoic acid,3-methyl-, phenylmethyl ester CAS No.:103-38-8 Molecular Structure: Molecular Structure of 103-38-8 (Butanoic acid,3-methyl-, phenylmethyl ester) Formula: C12H16 O2 Molecular Weight : 192.28 Synonyms: Isovalericacid, benzyl ester (6CI,8CI); Benzyl 3-methylbutanoate; Benzyl3-methylbutyrate; Benzyl

1-Methyl-L-4,5-dihydroorotic acid CAS 103365-69-1

Name: 1-Methyl-L-4,5-dihydroorotic acid CAS No.:103365-69-1 Molecular Structure: Molecular Structure of 103365-69-1 (1-Methyl-L-4,5-dihydroorotic acid) Formula: C6H8N2O4 Molecular Weight : 172.14 Synonyms: 4-Pyrimidinecarboxylicacid, hexahydro-1-methyl-2,6-dioxo-,(

(1R)-(-)-CAMPHORQUINONE CAS 10334-26-6

Name: (1R,4S)-1,7,7-Trimethylbicyclo[2.2.1]heptane-2,3-dione CAS No.:10334-26-6 Molecular Structure: Molecular Structure of 10334-26-6 ((1R,4S)-1,7,7-Trimethylbicyclo[2.2.1]heptane-2,3-dione) Formula: C10H14O2 Molecular Weight : 166.22 Synonyms: 2,3-Bornanedione,(57264285,1R,4S)-(-)-

3-Oxo-4-aza-5-alpha-androstane-17-beta-carboxylic acid CAS 103335-55-3

Name: 3-Oxo-4-aza-5-alpha-androstane-17-beta-carboxylic acid CAS No.:103335-55-3 Molecular Structure: Molecular Structure of 103335-55-3 (3-Oxo-4-aza-5-alpha-androstane-17-beta-carboxylic acid) Formula: C19H29NO3 Molecular Weight : 319.44 Synonyms: 4-Azaandrostane-17-carboxylicacid, 3-oxo-,(

Potassium saccharate CAS 10332-51-1

Name: 1,2-Benzisothiazol-3(2H)-one,1,1-dioxide, potassium salt (1:1) CAS No.:10332-51-1 Molecular Structure: Molecular Structure of 10332-51-1 (1,2-Benzisothiazol-3(2H)-one,1,1-dioxide, potassium salt (1:1)) Formula: C7H5 N O3 S . K Molecular Weight : 286.31952 Synonyms: 1,2-Benzisothiazol-3(2H)-one,1,1-dioxide,

Sodium perborate monohydrate CAS 10332-33-9

Name: Perboric acid(HBO(O2)), sodium salt, monohydrate (9CI) CAS No.:10332-33-9 Molecular Structure: Molecular Structure of 10332-33-9 (Perboric acid(HBO(O2)), sodium salt, monohydrate (9CI)) Formula: NaBO3.H2O Molecular Weight : 99.81 Synonyms: Perboricacid (HBO3), sodium salt, monohydrate (8CI);Interox A;Sodium

4-(pyridin-3-yl)cyclohexanone CAS 103319-09-1

Name: Cyclohexanone, 4-(3-pyridinyl)- CAS No.:103319-09-1 Molecular Structure: Molecular Structure of 103319-09-1 (Cyclohexanone, 4-(3-pyridinyl)-) Formula: C11H13NO Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport

N,N'-Bis(3-methylphenyl)-N,N'-diphenyl-9,9-spirobifluorene-2,7-diamine CAS 1033035-83-4

Name: 9,9'-Spirobi[9H-fluorene]-2',7'-diamine,N2',N7'-bis(3-methylphenyl)-N2',N7'-diphenyl- CAS No.:1033035-83-4 Molecular Structure: Molecular Structure of 1033035-83-4 (9,9'-Spirobi[9H-fluorene]-2',7'-diamine,N2',N7'-bis(3-methylphenyl)-N2',N7'-diphenyl-) Formula: C51H38N2 Molecular Weight : 678.86 Synonyms: LumTec

Ceritinib (LDK378) CAS 1032900-25-6

Name: LDK-378 CAS No.:1032900-25-6 Molecular Structure: Molecular Structure of 1032900-25-6 (LDK-378) Formula: Molecular Weight : Synonyms: Ceritinib; LDK 378; 5-chloro-n4-(2-((1-methylethyl)sulfonyl)phenyl)-n2-(5-methyl-2-(1-methylethoxy)-4-(4-piperidinyl)phenyl)-2,4-pyrimidinediamine EINECS: Density: Melting Point:

(S)-1-[4-[[2-(2-Aminopyrimidin-5-yl)-7-methyl-4-(morpholin-4-yl)thieno[3,2-d]pyrimidin-6-yl]methyl]piperazin-1-yl]-2-hydroxypropan-1-one CAS 1032754-93-0

Name: (2S)-1-(4-{[2-(2-Amino-5-pyrimidinyl)-7-methyl-4-(4-morpholinyl)thieno[3,2-d]pyrimidin-6-yl]methyl}-1-piperazinyl)-2-hydroxy-1-propanone CAS No.:1032754-93-0 Molecular Structure: Molecular Structure of 1032754-93-0

AZELAIC ACID DI(2-ETHYLHEXYL) ESTER CAS 103-24-2

Name: Nonanedioic acid,1,9-bis(2-ethylhexyl) ester CAS No.:103-24-2 Molecular Structure: Molecular Structure of 103-24-2 (Nonanedioic acid,1,9-bis(2-ethylhexyl) ester) Formula: C25H48 O4 Molecular Weight : 412.73 Synonyms: Azelaicacid, bis(2-ethylhexyl) ester (6CI,7CI,8CI);Nonanedioic acid,bis(2-ethylhexyl) ester

beta-D-(-)-Arabinose CAS 10323-20-3

Name: D-(-)-Arabinose CAS No.:10323-20-3 Molecular Structure: Molecular Structure of 10323-20-3 (D-(-)-Arabinose) Formula: C5H10O5 Molecular Weight : 150.13 Synonyms: Arabinose,D- (8CI);(-)-Arabinose;Anhydroarabinose;D-Arabinose; EINECS: 233-708-5 Density: 1.757 g/cm3 Melting Point: Boiling Point: 415.462 °C at 760

5-bromo-2-cyanobenzoic acid CAS 1032231-28-9

Name: 5-Bromo-2-cyanobenzoic acid CAS No.:1032231-28-9 Molecular Structure: Molecular Structure of 1032231-28-9 (5-Bromo-2-cyanobenzoic acid) Formula: Molecular Weight : 226 Synonyms: 5-bromo-2-cyanobenzoic acid EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk

4-(2-Chlorophenoxy)-N-[3-[(methylamino)carbonyl]phenyl]-1-Piperidinecarboxamide CAS 1032229-33-6

Name: 1-Piperidinecarboxamide,4-(2-chlorophenoxy)-N-[3-[(methylamino)carbonyl]phenyl]- CAS No.:1032229-33-6 Molecular Structure: Molecular Structure of 1032229-33-6 (1-Piperidinecarboxamide,4-(2-chlorophenoxy)-N-[3-[(methylamino)carbonyl]phenyl]-) Formula: C20H22ClN3O3 Molecular Weight : 387.86 Synonyms:

VAPREOTIDE CAS 103222-11-3

Name: L-Tryptophanamide,D-phenylalanyl-L-cysteinyl-L-tyrosyl-D-tryptophyl-L-lysyl-L-valyl-L-cysteinyl-,cyclic (2®7)-disulfide CAS No.:103222-11-3 Molecular Structure: Molecular Structure of 103222-11-3 (L-Tryptophanamide,D-phenylalanyl-L-cysteinyl-L-tyrosyl-D-tryptophyl-L-lysyl-L-valyl-L-cysteinyl-,cyclic

N-trans-caffeoyltyramine CAS 103188-48-3

Name: 2-Propenamide, 3-(3,4-dihydroxyphenyl)-N-[2-(4-hydroxyphenyl)ethyl]-,(57264268,2E)- CAS No.:103188-48-3 Molecular Structure: Molecular Structure of 103188-48-3 (2-Propenamide, 3-(3,4-dihydroxyphenyl)-N-[2-(4-hydroxyphenyl)ethyl]-,(57264269,2E)-) Formula: C17H17NO4 Molecular Weight : Synonyms: EINECS: Density:

(S)-Amlodipine CAS 103129-82-4

Name: (S)-Amlodipine CAS No.:103129-82-4 Molecular Structure: Molecular Structure of 103129-82-4 ((S)-Amlodipine) Formula: C20H25ClN2O5 Molecular Weight : 408.88 Synonyms: 3,5-Pyridinedicarboxylicacid, 2-[(2-aminoethoxy)methyl]-4-(2-chlorophenyl)-1,4-dihydro-6-methyl-,3-ethyl 5-methyl ester,(

2-AMINOTEREPHTHALIC ACID CAS 10312-55-7

Name: 2-Aminoterephthalic acid CAS No.:10312-55-7 Molecular Structure: Molecular Structure of 10312-55-7 (2-Aminoterephthalic acid) Formula: C8H7NO4 Molecular Weight : 181.15 Synonyms: Aminoterephthalic acid; EINECS: Density: 1.551 g/cm3 Melting Point: Boiling Point: 450.7 °C at 760 mmHg Flash Point: 226.4 °C

Methyl triphenyl phosphonium chloride CAS 1031-15-8

Name: Phosphonium,methyltriphenyl-, chloride (1:1) CAS No.:1031-15-8 Molecular Structure: Molecular Structure of 1031-15-8 (Phosphonium,methyltriphenyl-, chloride (1:1)) Formula: C19H18ClP Molecular Weight : 312.78 Synonyms: Methyltriphenylphosphoniumchloride (6CI,7CI);Phosphonium, methyltriphenyl-, chloride

3-Ethyl-5-methyl-4-(2-chlorophenyl)-2-(2,2-diethoxy-ethoxymethyl)-6-methyl-1,4-dihydropyridine-3,5-dicarboxylate CAS 103094-30-0

Name: 3-Ethyl-5-methyl-4-(2-chlorophenyl)-2-(2,2-diethoxy-ethoxymethyl)-6-methyl-1,4-dihydropyridine-3,5-dicarboxylate CAS No.:103094-30-0 Molecular Structure: Molecular Structure of 103094-30-0 (3-Ethyl-5-methyl-4-(2-chlorophenyl)-2-(2,2-diethoxy-ethoxymethyl)-6-methyl-1,4-dihydropyridine-3,5-dicarboxylate) Formula:

2-Ethylhexyl acetate CAS 103-09-3

Name: Acetic acid,2-ethylhexyl ester CAS No.:103-09-3 Molecular Structure: Molecular Structure of 103-09-3 (Acetic acid,2-ethylhexyl ester) Formula: C10H20O2 Molecular Weight : 172.30 Synonyms: b-Ethylhexyl acetate;Octyl acetate;OctASOLV;2-Ethylhexyl acetate;NSC 8897;1-Hexanol,2-ethyl-, acetate (6CI);2-Ethyl-1-hexanol

Aminoguanidinium nitrate CAS 10308-82-4

Name: Aminoguanidine nitrate CAS No.:10308-82-4 Molecular Structure: Molecular Structure of 10308-82-4 (Aminoguanidine nitrate) Formula: CH6N4.HNO3 Molecular Weight : 137.10 Synonyms: Monoaminoguanidium nitrate;Guanidine, amino-, mononitrate;Hydrazinecarboximidamide, mononitrate;Aminoguanidinenitrate;Aminoguanidine,

D-Glucopyranoside, Methyl 1-C-[3-[[5-(4-fluorophenyl)-2-thienyl]Methyl]-4-Methylphenyl]- CAS 1030825-21-8

Name: Methyl 1-C-(3-{[5-(4-fluorophenyl)-2-thienyl]methyl}-4-methylphenyl)-D-glucopyranoside CAS No.:1030825-21-8 Molecular Structure: Molecular Structure of 1030825-21-8 (Methyl 1-C-(3-{[5-(4-fluorophenyl)-2-thienyl]methyl}-4-methylphenyl)-D-glucopyranoside) Formula: C25H27FO6S Molecular Weight : 474.544183 Synonyms:

1-Bromo-3,5-diphenylbenzene CAS 103068-20-8

Name: 1-Bromo-3,5-diphenylbenzene CAS No.:103068-20-8 Molecular Structure: Molecular Structure of 103068-20-8 (1-Bromo-3,5-diphenylbenzene) Formula: C18H13Br Molecular Weight : 309.1998 Synonyms: 3,5-Diphenyl-1-bromophenyl EINECS: Density: 1.309 g/cm3 Melting Point: Boiling Point: 401.13 °C at 760 mmHg Flash Point:

2-METHYL-4-PHENYL-2-BUTANOL CAS 103-05-9

Name: Benzenepropanol, a,a-dimethyl- CAS No.:103-05-9 Molecular Structure: Molecular Structure of 103-05-9 (Benzenepropanol, a,a-dimethyl-) Formula: C11H16O Molecular Weight : 164.24 Synonyms: 2-Butanol,2-methyl-4-phenyl-

Lufenuron CAS 103055-07-8

Name: Lufenuron CAS No.:103055-07-8 Molecular Structure: Molecular Structure of 103055-07-8 (Lufenuron) Formula: C17H8Cl2F8N2O3 Molecular Weight : 511.15 Synonyms: Fluphenacur;1-(2,5-Dichloro-4-(1,1,2,3,3,3-hexafluoropropoxy)phenyl)-3-(2,6-difluorobenzoyl)urea;Benzamide,N-[[[2,5-dichloro-4-(1,1,2,3,3,3-

(PHENYLTHIO)ACETIC ACID CAS 103-04-8

Name: Acetic acid,2-(phenylthio)- CAS No.:103-04-8 Molecular Structure: Molecular Structure of 103-04-8 (Acetic acid,2-(phenylthio)-) Formula: C8H8 O2 S Molecular Weight : 152.2135 Synonyms: Aceticacid,(57264255,phenylthio)- (6CI,7CI,8CI,9CI); (Phenylthio)acetic acid;(S-Phenylmercapto)acetic acid;

4,6-Bis(3,5-di(pyridin-4-yl)phenyl)-2-MethylpyriMidine CAS 1030380-51-8

Name: 4,6-bis(3,5-di(pyridin-4-yl)phenyl)-2-methylpyrimidine CAS No.:1030380-51-8 Molecular Structure: Molecular Structure of 1030380-51-8 (4,6-bis(3,5-di(pyridin-4-yl)phenyl)-2-methylpyrimidine) Formula: C37H26N6 Molecular Weight : 554.64 Synonyms: B4PYMPM; EINECS: Density: Melting Point: Boiling Point: Flash Point:

4,6-bis(3,5-dichlorophenyl)-2-methylpyrimidine CAS 1030380-50-7

0.99

3,3'-(5-bromo-1,3-phenylene)dipyridine CAS 1030380-36-9

0.99

Anilinoacetic acid CAS 103-01-5

Name: Glycine,N-phenyl- CAS No.:103-01-5 Molecular Structure: Molecular Structure of 103-01-5 (Glycine,N-phenyl-) Formula: C8H9NO2 Molecular Weight : 151.16 Synonyms: H-DL-Phg-OH;TR-NPG;(Phenylamino)aceticacid;Acetic acid,(57264250,phenylamino)-;Anilinoacetic acid;N-(Phenylamino)aceticacid;N-Phenylglycine;NSC 83567;

3'-O-Methyl-D-adenosine CAS 10300-22-8

Name: Adenosine, 3'-O-methyl-(7CI,8CI,9CI) CAS No.:10300-22-8 Molecular Structure: Molecular Structure of 10300-22-8 (Adenosine, 3'-O-methyl-(7CI,8CI,9CI)) Formula: C11H15N5O4 Molecular Weight : 281.2679 Synonyms: 3'-O-Methyladenosine;NSC 103062; EINECS: Density: 1.84 g/cm3 Melting Point: Boiling Point: 623.8 °C at

8-AZAADENOSINE CAS 10299-44-2

Name: 3H-1,2,3-Triazolo[4,5-d]pyrimidin-7-amine,3-β-D-ribofuranosyl- CAS No.:10299-44-2 Molecular Structure: Molecular Structure of 10299-44-2 (3H-1,2,3-Triazolo[4,5-d]pyrimidin-7-amine,3-β-D-ribofuranosyl-) Formula: C9H12N6O4 Molecular Weight : 268.2294 Synonyms:

bullatacin CAS 102989-24-2

Name: 2(5H)-Furanone,3-[(2R,13R)-2,13-dihydroxy-13-[(2R,2'R,5R,5'R)-octahydro-5'-[(1R)-1-hydroxyundecyl][2,2'-bifuran]-5-yl]tridecyl]-5-methyl-,(57264245,5S)- CAS No.:102989-24-2 Molecular Structure: Molecular Structure of 102989-24-2

Plx-4032 (RG7024) CAS 1029872-54-5

Name: 1-Propanesulfonamide, N-[3-[[5-(4-chlorophenyl)-1H-pyrrolo[2,3-b]pyridin-3-yl]carbonyl]-2,4-difluorophenyl]- CAS No.:1029872-54-5 Molecular Structure: Molecular Structure of 1029872-54-5 (1Propanesulfonamide, N-[3-[[5-(4-chlorophenyl)-1H-pyrrolo[2,3-b]pyridin-3-yl]carbonyl]-2,4-difluorophenyl]-) Formula:

INCB28060 CAS 1029712-80-8

Name: Benzamide, 2-fluoro-N-methyl-4-[7-(6-quinolinylmethyl)imidazo[1,2-b][1,2,4]triazin-2-yl]- CAS No.:1029712-80-8 Molecular Structure: Molecular Structure of 1029712-80-8 (Benzamide, 2-fluoro-N-methyl-4-[7-(6-quinolinylmethyl)imidazo[1,2-b][1,2,4]triazin-2-yl]-) Formula: C23H17FN6O Molecular Weight : Synonyms:

Universal CAS 102962-70-9

Name: Universal CAS No.:102962-70-9 Molecular Structure: Molecular Structure of 102962-70-9 (Universal) Formula: Unspecified Molecular Weight : 0 Synonyms: Universal EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

Cesium sulfate CAS 10294-54-9

Name: Cesium sulfate CAS No.:10294-54-9 Molecular Structure: Molecular Structure of 10294-54-9 (Cesium sulfate) Formula: Cs2SO4 Molecular Weight : 361.87 Synonyms: Cesiumsulfate (Cs2(SO4)) (6CI,7CI);Sulfuric acid, dicesium salt (8CI,9CI);Cesiumsulfate;Dicesium sulfate;Sulfuricacid, cesium salt (1:2); EINECS: 233-662-6

CERIUM(IV) SULFATE TETRAHYDRATE CAS 10294-42-5

Name: Sulfuric acid,cerium(4+) salt (2:1), tetrahydrate (8CI,9CI) CAS No.:10294-42-5 Molecular Structure: Molecular Structure of 10294-42-5 (Sulfuric acid,cerium(4+) salt (2:1), tetrahydrate (8CI,9CI)) Formula: Ce.2H2O4S.4H2O Molecular Weight : 404.29 Synonyms: Cericsulfate tetrahydrate;Cerium disulfate

CERIUM(III) NITRATE HEXAHYDRATE CAS 10294-41-4

Name: Nitric acid, cerium(3+)salt, hexahydrate (8CI,9CI) CAS No.:10294-41-4 Molecular Structure: Molecular Structure of 10294-41-4 (Nitric acid, cerium(3+)salt, hexahydrate (8CI,9CI)) Formula: Ce(NO3)3.6(H2O) Molecular Weight : 434.22 Synonyms: Cerium(III) nitrate hexahydrate;Cerium nitrate hexahydrate;Cerium

Barium chromate CAS 10294-40-3

Name: Chromic acid (H2CrO4),barium salt (1:1) CAS No.:10294-40-3 Molecular Structure: Molecular Structure of 10294-40-3 (Chromic acid (H2CrO4),barium salt (1:1)) Formula: BaCrO4 Molecular Weight : 253.32 Synonyms: Bariumchromate(VI) (6CI);Barium chromate (1:1);Barium chromate(BaCrO4);Barium chromium oxide

Boron trichloride CAS 10294-34-5

Name: Borane, trichloro- CAS No.:10294-34-5 Molecular Structure: Molecular Structure of 10294-34-5 (Borane, trichloro-) Formula: BCl3 Molecular Weight : 117.16 Synonyms: Boronchloride (BCl3) (8CI);Trichloroborane;Trichloroboron; EINECS: 233-658-4 Density: 1.384 g/cm3 Melting Point: Boiling Point: 12.5 °C at 760 mmHg

GOLD CHLORIDE CAS 10294-30-1

Name: Gold chloride (AuCl3),dihydrate (8CI,9CI) CAS No.:10294-30-1 Molecular Structure: Molecular Structure of 10294-30-1 (Gold chloride (AuCl3),dihydrate (8CI,9CI)) Formula: AuCl3 . 2 H2 O Molecular Weight : 339.79 Synonyms: Auricchloride dihydrate EINECS: 240-948-4 Density: 3.9g/mLat25℃ Melting Point: Boiling Point:

Silver sulfate CAS 10294-26-5

Name: Sulfuric acid,silver(1+) salt (1:2) CAS No.:10294-26-5 Molecular Structure: Molecular Structure of 10294-26-5 (Sulfuric acid,silver(1+) salt (1:2)) Formula: Ag2SO4 Molecular Weight : 311.79 Synonyms: Disilver monosulfate;Disilver(1+) sulfate;Silver sulfate (Ag2SO4);Sulfuric acid silver salt (1:2); EINECS:

99% CAS 102939-46-8

Name: Hexanoic acid, 6-chloro-6-oxo- CAS No.:102939-46-8 Molecular Structure: Molecular Structure of 102939-46-8 (Hexanoic acid, 6-chloro-6-oxo-) Formula: C6H9ClO3 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport

Cinnamoyl chloride CAS 102-92-1

Name: 2-Propenoyl chloride,3-phenyl- CAS No.:102-92-1 Molecular Structure: Molecular Structure of 102-92-1 (2-Propenoyl chloride,3-phenyl-) Formula: C9H7ClO Molecular Weight : 166.61 Synonyms: Cinnamoylchloride (6CI,7CI,8CI);3-Phenyl-2-propenoyl chloride;3-Phenylacryloylchloride;3-Phenylpropenoyl

3-(4-bromophenyl)-N-phenylcarbazole CAS 1028647-93-9

Name: 3-(4-Bromophenyl)-9-phenyl-9H-Carbazole CAS No.:1028647-93-9 Molecular Structure: Molecular Structure of 1028647-93-9 (3-(4-Bromophenyl)-9-phenyl-9H-Carbazole) Formula: C24H16BrN Molecular Weight : 398.29 Synonyms: 3-(4-bromophenyl)-N-phenyl-9H-Carbazole; EINECS: Density: 1.329 g/cm3 Melting Point: Boiling

(2S)-2-(2-Oxopyrrolidin-1-yl)butanoic acid CAS 102849-49-0

Name: (2S)-2-(2-Oxopyrrolidin-1-yl)butanoic acid CAS No.:102849-49-0 Molecular Structure: Molecular Structure of 102849-49-0 ((2S)-2-(2-Oxopyrrolidin-1-yl)butanoic acid) Formula: C8H13NO3 Molecular Weight : 171.20 Synonyms: UCB-L 057;1-Pyrrolidineaceticacid, a-ethyl-2-oxo-,(57264230,S)-;2-Pyrrolidinone-n-butyric

D-Pinitol CAS 10284-63-6

Name: D-chiro-Inositol,3-O-methyl- CAS No.:10284-63-6 Molecular Structure: Molecular Structure of 10284-63-6 (D-chiro-Inositol,3-O-methyl-) Formula: C7H14O6 Molecular Weight : 194.18 Synonyms: Inositol,3-O-methyl-, D-chiro- (8CI);Pinitol,(57264228,+)-

TRIALLYL PHOSPHITE CAS 102-84-1

Name: Triallyl phosphite CAS No.:102-84-1 Molecular Structure: Molecular Structure of 102-84-1 (Triallyl phosphite) Formula: C9H15 O3 P Molecular Weight : 202.21 Synonyms: Allylphosphite (6CI,7CI); Phosphorous acid, tri-2-propenyl ester (9CI); Phosphorousacid, triallyl ester (8CI); NSC 44606; Triallyl phosphite;

2-CHLORO-3,4-DIMETHYLPHENOL CAS 10283-15-5

Name: Phenol,2-chloro-3,4-dimethyl- CAS No.:10283-15-5 Molecular Structure: Molecular Structure of 10283-15-5 (Phenol,2-chloro-3,4-dimethyl-) Formula: C8H9 Cl O Molecular Weight : 156.6095 Synonyms: 3,4-Xylenol,2-chloro- (7CI,8CI); 2-Chloro-3,4-dimethylphenol; 2-Chloro-3,4-xylenol;3,4-Dimethyl-2-chlorophenol; NSC

N,N-DIBUTYL-1,3-PROPANEDIAMINE CAS 102-83-0

Name: 1,3-Propanediamine,N1,N1-dibutyl- CAS No.:102-83-0 Molecular Structure: Molecular Structure of 102-83-0 (1,3-Propanediamine,N1,N1-dibutyl-) Formula: C11H26N2 Molecular Weight : 186.34 Synonyms: 1,3-Propanediamine,N,N-dibutyl-

CyclopropanecarboxaMide, 1-aMino-N-(cyclopropylsulfonyl)-2-ethenyl-,(57264223,1R,2S)-, 4-Methylbenzenesulfonate(1:1) CAS 1028252-16-5

0.99

2,2'-(Butylimino)diethanol CAS 102-79-4

Name: Ethanol,2,2'-(butylimino)bis- CAS No.:102-79-4 Molecular Structure: Molecular Structure of 102-79-4 (Ethanol,2,2'-(butylimino)bis-) Formula: C8H19NO2 Molecular Weight : 161.24 Synonyms: Ethanol,2,2'-(butylimino)di-

2-(Morpholinothio)benzothiazole CAS 102782-92-3

Name: Siloxanes andSilicones, 3-[(2-aminoethyl)amino]propyl Me, di-Me, methoxy-terminated CAS No.:102782-92-3 Molecular Structure: Molecular Structure of 102782-92-3 (Siloxanes andSilicones, 3-[(2-aminoethyl)amino]propyl Me, di-Me, methoxy-terminated) Formula: Molecular Weight : 0 Synonyms: Siloxanes and Silicones,

2-(Morpholinothio)benzothiazole CAS 102-77-2

Name: 2-Benzothiazolesulfenemorpholide CAS No.:102-77-2 Molecular Structure: Molecular Structure of 102-77-2 (2-Benzothiazolesulfenemorpholide) Formula: C11H12N2OS2 Molecular Weight : 252.35 Synonyms: Benzothiazole,2-(morpholinothio)- (6CI,7CI,8CI);Morpholine, 4-(2-benzothiazolylthio)-

4,4,5A-TRIMETHYLPERHYDRO-1-BENZOXIREN-2-ONE CAS 10276-21-8

Name: 7-Oxabicyclo[4.1.0]heptan-2-one,4,4,6-trimethyl- CAS No.:10276-21-8 Molecular Structure: Molecular Structure of 10276-21-8 (7-Oxabicyclo[4.1.0]heptan-2-one,4,4,6-trimethyl-) Formula: C9H14O2 Molecular Weight : 154.21 Synonyms:

99% CAS 1027524-44-2

0.99

2-(diethylamino)-N-(2,4,6-trimethylphenyl)acetamide monohydrochloride CAS 1027-14-1

Name: Acetamide,2-(diethylamino)-N-(2,4,6-trimethylphenyl)-, hydrochloride (1:1) CAS No.:1027-14-1 Molecular Structure: Molecular Structure of 1027-14-1 (Acetamide,2-(diethylamino)-N-(2,4,6-trimethylphenyl)-, hydrochloride (1:1)) Formula: C15H24 N2 O . Cl H Molecular Weight : 284.87 Synonyms:

TEREPHTHALIC ACID DIALLYL ESTER CAS 1026-92-2

Name: 1,4-Benzenedicarboxylicacid, 1,4-di-2-propen-1-yl ester CAS No.:1026-92-2 Molecular Structure: Molecular Structure of 1026-92-2 (1,4-Benzenedicarboxylicacid, 1,4-di-2-propen-1-yl ester) Formula: C14H14O4 Molecular Weight : 246.26 Synonyms: 1,4-Benzenedicarboxylicacid, di-2-propenyl ester (9CI);Terephthalic acid,

2-Cyanoethyl N,N,N',N'-tetraisopropylphosphorodiamidite CAS 102691-36-1

Name: Phosphorodiamidousacid, N,N,N',N'-tetrakis(1-methylethyl)-, 2-cyanoethyl ester CAS No.:102691-36-1 Molecular Structure: Molecular Structure of 102691-36-1 (Phosphorodiamidousacid, N,N,N',N'-tetrakis(1-methylethyl)-, 2-cyanoethyl ester) Formula: C15H32N3OP Molecular Weight : 301.41 Synonyms:

Cyclopentanone,2,5-bis(2-furanylmethylene)- CAS 1026-78-4

Name: Cyclopentanone,2,5-bis(2-furanylmethylene)- CAS No.:1026-78-4 Molecular Structure: Molecular Structure of 1026-78-4 (Cyclopentanone,2,5-bis(2-furanylmethylene)-) Formula: C15H12 O3 Molecular Weight : 240.254 Synonyms: Cyclopentanone,2,5-difurfurylidene- (6CI,7CI,8CI);

LB-100 CAS 1026680-07-8

Name: LB-100 CAS No.:1026680-07-8 Molecular Structure: Molecular Structure of 1026680-07-8 (LB-100) Formula: Molecular Weight : 0 Synonyms: LB-100;3-(1-methylpiperazine-4-carbonyl)-7-oxa-bicyclo[2.2.1]heptane-2-carboxylic acid;LB-100 3-(4-METHYLPIPERAZINE-1-CARBONYL)-7-OXABICYCLO[2.2.1]HEPTANE-2-CARBOXYLIC ACID;methyl

2-Bromoacetyl-6-methoxynaphtalene CAS 10262-65-4

Name: Ethanone, 2-bromo-1-(6-methoxy-2-naphthalenyl)- CAS No.:10262-65-4 Molecular Structure: Molecular Structure of 10262-65-4 (Ethanone, 2-bromo-1-(6-methoxy-2-naphthalenyl)-) Formula: C13H11BrO2 Molecular Weight : 279.12924 Synonyms:

3-Cyclohexylamino-2-hydroxypropanesulfonic acid sodium salt CAS 102601-34-3

Name: 1-Propanesulfonic acid,3-(cyclohexylamino)-2-hydroxy-, sodium salt (1:1) CAS No.:102601-34-3 Molecular Structure: Molecular Structure of 102601-34-3 (1-Propanesulfonic acid,3-(cyclohexylamino)-2-hydroxy-, sodium salt (1:1)) Formula: C9H18NNaO4S Molecular Weight : 259.30 Synonyms: 1-Propanesulfonicacid,

2,7-Dibromopyrene CAS 102587-98-4

Name: 2,7-Dibromopyrene CAS No.:102587-98-4 Molecular Structure: Molecular Structure of 102587-98-4 (2,7-Dibromopyrene) Formula: C16H8Br2 Molecular Weight : 357.89 Synonyms: 2,7-Dibromopyrene EINECS: Density: 1.853 g/cm3 Melting Point: Boiling Point: 469.56 °C at 760 mmHg Flash Point: 277.277 °C Solubility:

99% CAS 10258-02-3

Name: 4(1H)-Quinolinone, 2,3-dihydro-1-phenyl-, oxime CAS No.:10258-02-3 Molecular Structure: Molecular Structure of 10258-02-3 (4(1H)-Quinolinone, 2,3-dihydro-1-phenyl-, oxime) Formula: C15H14N2O Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard

Calcium sulfite CAS 10257-55-3

Name: Sulfurous acid, calciumsalt (1:1) CAS No.:10257-55-3 Molecular Structure: Molecular Structure of 10257-55-3 (Sulfurous acid, calciumsalt (1:1)) Formula: Ca. H2 O3 S Molecular Weight : 136.1406 Synonyms: Calciumsulfite (6CI); Calcium sulfite (CaSO3) (7CI); Calcium sulfite (1:1) EINECS: Density: g/cm3 Melting

- ## Copper sulfate monohydrate CAS 10257-54-2

  Name: Sulfuric acidcopper(2+) salt (1:1), monohydrate (8CI,9CI) CAS No.:10257-54-2 Molecular Structure: Molecular Structure of 10257-54-2 (Sulfuric acidcopper(2+) salt (1:1), monohydrate (8CI,9CI)) Formula: CuSO4.H2O Molecular Weight : 177.62 Synonyms: Coppermesosulfate (Cu(SO4).H2O) (7CI);Copper sulfate (CuSO4)

- ## 2,5-Dimethoxyaniline CAS 102-56-7

  Name: Benzenamine,2,5-dimethoxy- CAS No.:102-56-7 Molecular Structure: Molecular Structure of 102-56-7 (Benzenamine,2,5-dimethoxy-) Formula: C8H11NO2 Molecular Weight : 153.18 Synonyms: Aniline,2,5-dimethoxy- (7CI,8CI);1-Amino-2,5-dimethoxybenzene;2,5-Dimethoxybenzenamine;Aminohydroquinone dimethyl ether;C.I.

- ## MERCAPTOPYRUVIC ACID SODIUM SALT CAS 10255-67-1

  Name: Propanoic acid,3-mercapto-2-oxo-, sodium salt (1:1) CAS No.:10255-67-1 Molecular Structure: Molecular Structure of 10255-67-1 (Propanoic acid,3-mercapto-2-oxo-, sodium salt (1:1)) Formula: C3H4 O3 S . Na Molecular Weight : 142.1089 Synonyms: Propanoicacid, 3-mercapto-2-oxo-, monosodium salt (9CI); Pyruvic acid,

- ## NORSOLORINICACID CAS 10254-99-6

  Name: 9,10-Anthracenedione,1,3,6,8-tetrahydroxy-2-(1-oxohexyl)- CAS No.:10254-99-6 Molecular Structure: Molecular Structure of 10254-99-6 (9,10-Anthracenedione,1,3,6,8-tetrahydroxy-2-(1-oxohexyl)-) Formula: C20H18 O7 Molecular Weight : 0 Synonyms: Anthraquinone,2-hexanoyl-1,3,6,8-tetrahydroxy- (7CI,8CI);

- ## 3,4-Dichlorobenzylamine CAS 102-49-8

  Name: Benzenemethanamine,3,4-dichloro- CAS No.:102-49-8 Molecular Structure: Molecular Structure of 102-49-8 (Benzenemethanamine,3,4-dichloro-) Formula: C7H7Cl2N Molecular Weight : 176.04 Synonyms: Benzylamine,3,4-dichloro-

- ## 1H,1H,2H,2H-PERFLUOROOCTYLDIMETHYLCHLOROSILANE CAS 102488-47-1

  Name: Silane,chlorodimethyl(3,3,4,4,5,5,6,6,7,7,8,8,8-tridecafluorooctyl)- CAS No.:102488-47-1 Molecular Structure: Molecular Structure of 102488-47-1 (Silane,chlorodimethyl(3,3,4,4,5,5,6,6,7,7,8,8,8-tridecafluorooctyl)-) Formula: C10H10ClF13Si Molecular Weight : 440.00 Synonyms:

- ## 1,2-Dichloro-4-(chloromethyl)benzene CAS 102-47-6

  Name: Benzene,1,2-dichloro-4-(chloromethyl)- CAS No.:102-47-6 Molecular Structure: Molecular Structure of 102-47-6 (Benzene,1,2-dichloro-4-(chloromethyl)-) Formula: C7H5Cl3 Molecular Weight : 195.47 Synonyms: Toluene,a,3,4-trichloro-

- ## 11,12-Dihydro-11-phenylindolo[2,3-a]carbazole CAS 1024598-06-8

  Name: 11-phenylindolo[2,3-a]carbazole CAS No.:1024598-06-8 Molecular Structure: Molecular Structure of 1024598-06-8 (11-phenylindolo[2,3-a]carbazole) Formula: Molecular Weight : 332.40444 Synonyms: 11,12-dihydro-11-phenyl-Indolo[2,3-a]carbazole; Indolo[2,3-a]carbazole, 11,12-dihydro-11-phenyl-; EINECS: Density:

- ## 6-NITRO-2-OXO-2H-CHROMENE-3-CARBOXYLIC ACID CAS 10242-15-6

  Name: 2H-1-Benzopyran-3-carboxylicacid, 6-nitro-2-oxo- CAS No.:10242-15-6 Molecular Structure: Molecular Structure of 10242-15-6 (2H-1-Benzopyran-3-carboxylicacid, 6-nitro-2-oxo-) Formula: C10H5 N O6 Molecular Weight : 235.15 Synonyms: 6-Nitrocoumarin-3-carboxylicacid; NSC 74856 EINECS: Density: 1.689g/cm3 Melting

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99% CAS 102390-86-3
0.99

- ## 5-CARBOXYOXINDOLE CAS 102359-00-2
  Name: 5-Carboxyoxindole CAS No.:102359-00-2 Molecular Structure: Molecular Structure of 102359-00-2 (5-Carboxyoxindole) Formula: C9H7NO3 Molecular Weight : 177.1568 Synonyms: 5-CARBOXYOXINDOLE;RARECHEM AL BO 0983;OXINDOLE-5-CARBOXYLIC ACID;2-Oxo-indoline-5-carboxylic acid;5-CARBOXY-2-OXINDOLE;5-Carboxyoxindole 97%

- ## omeprazole cyclic sulfenamide CAS 102332-89-8
  Name: 5H-Pyrido[1',2':4,5][1,2,4]thiadiazino[2,3-a]benzimidazol-13-ium(9CI) CAS No.:102332-89-8 Molecular Structure: Molecular Structure of 102332-89-8 (5H-Pyrido[1',2':4,5][1,2,4]thiadiazino[2,3-a]benzimidazol-13-ium(9CI)) Formula: C13H10 N3 S Molecular Weight : 0 Synonyms: omeprazole cyclic sulfenamide EINECS:

- ## ISOPROPYL LAURATE CAS 10233-13-3
  Name: Dodecanoic acid,1-methylethyl ester CAS No.:10233-13-3 Molecular Structure: Molecular Structure of 10233-13-3 (Dodecanoic acid,1-methylethyl ester) Formula: C15H30O2 Molecular Weight : 242.40 Synonyms: Lauricacid, isopropyl ester (6CI,7CI,8CI);Dodecanoic acid isopropyl ester;Emcol-IL;Isopropyl dodecanoate;

- ## DOPAC CAS 102-32-9
  Name: Benzeneacetic acid,3,4-dihydroxy- CAS No.:102-32-9 Molecular Structure: Molecular Structure of 102-32-9 (Benzeneacetic acid,3,4-dihydroxy-) Formula: C8H8O4 Molecular Weight : 168.16 Synonyms: Aceticacid,(57264192,3,4-dihydroxyphenyl)- (8CI);(3,4-Dihydroxyphenyl)acetic acid;3,4-Dihydroxybenzeneacetic acid;Ba

- ## 5-AMINO-1-N-METHYLINDOLE CAS 102308-97-4
  Name: 1H-Indol-5-amine,1-methyl- CAS No.:102308-97-4 Molecular Structure: Molecular Structure of 102308-97-4 (1H-Indol-5-amine,1-methyl-) Formula: C9H10N2 Molecular Weight : 146.19 Synonyms: Indole,5-amino-1-methyl-

- ## (S)-(-)-3-TERT-BUTOXYCARBONYL-4-FORMYL-2,2-DIMETHYL-1,3-OXAZOLIDINE CAS 102308-32-7
  Name: (S)-(-)-3-Boc-2,2-dimethyloxazolidine-4-carboxaldehyde CAS No.:102308-32-7 Molecular Structure: Molecular Structure of 102308-32-7 ((S)-(-)-3-Boc-2,2-dimethyloxazolidine-4-carboxaldehyde) Formula: C11H19NO4 Molecular Weight : 229.27 Synonyms: 3-Oxazolidinecarboxylicacid, 4-formyl-2,2-dimethyl-, 1,1-dimethylethyl

- ## 2,4-Dihydroxy-2,5-dimethyl-3(2H)-furan-3-one CAS 10230-62-3
  Name: 3(2H)-Furanone,2,4-dihydroxy-2,5-dimethyl- CAS No.:10230-62-3 Molecular Structure: Molecular Structure of 10230-62-3 (3(2H)-Furanone,2,4-dihydroxy-2,5-dimethyl-) Formula: C6H8 O4 Molecular Weight : 0 Synonyms: 2,5-Dimethyl-2,4-dihydroxy-3(2H)-furanone EINECS: Density: Melting Point: Boiling Point: Flash Point:

- ## Ethanone, 2-bromo-1-(2,2-dimethyl-4H-1,3-benzodioxin-6-yl)- CAS 102293-80-1
  Name: Ethanone, 2-bromo-1-(2,2-dimethyl-4H-1,3-benzodioxin-6-yl)- CAS No.:102293-80-1 Molecular Structure: Molecular Structure of 102293-80-1 (Ethanone, 2-bromo-1-(2,2-dimethyl-4H-1,3-benzodioxin-6-yl)-) Formula: C12H13BrO3 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point:

- ## Ethyl pyrazolo[1,5-a]pyriMidine-6-carboxylate CAS 1022920-59-7
  Name: Ethyl pyrazolo[1,5-a]pyrimidine-6-carboxylate CAS No.:1022920-59-7 Molecular Structure: Molecular Structure of 1022920-59-7 (Ethyl pyrazolo[1,5-a]pyrimidine-6-carboxylate) Formula: C9H9N3O2 Molecular Weight : 191.18666 Synonyms: Ethyl pyrazolo[1,5-a]pyriMidine-6-carboxylate;Pyrazolo[1,5-a]pyrimidine-6-carboxylic

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N1-(3-Aminophenyl)acetamide CAS 102-28-3

Name: Acetamide,N-(3-aminophenyl)- CAS No.:102-28-3 Molecular Structure: Molecular Structure of 102-28-3 (Acetamide,N-(3-aminophenyl)-) Formula: C8H10N2O Molecular Weight : 150.20 Synonyms: m-Aminoacetanilide;Acetanilide,3'-amino-

- ## 1-Bromo-5-hexanone CAS 10226-29-6

  Name: 1-Bromo-5-hexanone CAS No.:10226-29-6 Molecular Structure: Molecular Structure of 10226-29-6 (1-Bromo-5-hexanone) Formula: C6H11BrO Molecular Weight : 179.06 Synonyms: 1-Bromo-5-hexanone;5-Oxohexyl bromide;6-Bromo-2-hexanone; EINECS: 233-544-4 Density: 1.298 g/cm3 Melting Point: Boiling Point: 227.2 °C at 760

- ## 2,2-Dibromo-2-cyanoacetamide CAS 10222-01-2

  Name: 2,2-Dibromo-2-cyanoacetamide CAS No.:10222-01-2 Molecular Structure: Molecular Structure of 10222-01-2 (2,2-Dibromo-2-cyanoacetamide) Formula: C3H2Br2N2O Molecular Weight : 241.87 Synonyms: 2,2-Dibromo-3-nitrilopropionamide;2,2-Dibromo-3-nitrilopropanamide; EINECS: 233-539-7 Density: 2.451 g/cm3 Melting Point:

- ## 5-BROMO-2'-DEOXYURIDINE 5'-TRIPHOSPHATE SODIUM SALT CAS 102212-99-7

  Name: Uridine5'-(tetrahydrogen triphosphate), 5-bromo-2'-deoxy-, tetrasodium salt (9CI) CAS No.:102212-99-7 Molecular Structure: Molecular Structure of 102212-99-7 (Uridine5'-(tetrahydrogen triphosphate), 5-bromo-2'-deoxy-, tetrasodium salt (9CI)) Formula: C9H14 Br N2 O14 P3 . 4 Na Molecular Weight : 547.04 Synonyms:

- ## Phenethyl phenylacetate CAS 102-20-5

  Name: Benzeneacetic acid,2-phenylethyl ester CAS No.:102-20-5 Molecular Structure: Molecular Structure of 102-20-5 (Benzeneacetic acid,2-phenylethyl ester) Formula: C16H16O2 Molecular Weight : 240.32 Synonyms: Benzylcarbinyl a-toluate;NSC 6676;Phenethyl phenylacetate;Phenylethyl phenylacetate;b-Phenylethyl

- ## 1-(4-NITROPHENYL)PYRROLIDINE CAS 10220-22-1

  Name: Pyrrolidine,1-(4-nitrophenyl)- CAS No.:10220-22-1 Molecular Structure: Molecular Structure of 10220-22-1 (Pyrrolidine,1-(4-nitrophenyl)-) Formula: C10H12N2O2 Molecular Weight : 192.21 Synonyms: Pyrrolidine,1-(p-nitrophenyl)-

- ## 3-(pyridin-3-ylmethylamino)-1-(2,6,6-trimethylcyclohex-3-en-1-yl)butan-1-one CAS 1022-01-2

  0.99

- ## (2R)-α,α,8α,8aα-Tetramethyl-1,2,3,4,6,7,8,8a-octahydronaphthalene-2α-methanol CAS 10219-71-3

  Name: 2-Naphthalenemethanol,1,2,3,4,6,7,8,8a-octahydro-a,a,8,8a-tetramethyl-,(57264175,2R,8S,8aR)- CAS No.:10219-71-3 Molecular Structure: Molecular Structure of 10219-71-3 (2-Naphthalenemethanol,1,2,3,4,6,7,8,8a-octahydro-a,a,8,8a-tetramethyl-,(57264176,2R,8S,8aR)-) Formula: C15H26 O Molecular Weight : 0 Synonyms:

- ## N-Boc-cis-4-Hydroxy-L-proline methyl ester CAS 102195-79-9

  Name: 1,2-Pyrrolidinedicarboxylicacid, 4-hydroxy-, 1-(1,1-dimethylethyl) 2-methyl ester,(57264171,2S,4S)- CAS No.:102195-79-9 Molecular Structure: Molecular Structure of 102195-79-9 (1,2-Pyrrolidinedicarboxylicacid, 4-hydroxy-, 1-(1,1-dimethylethyl) 2-methyl ester,(57264172,2S,4S)-) Formula: C11H19NO5 Molecular

- ## BENZYL PHENYLACETATE CAS 102-16-9

  Name: Benzeneacetic acid,phenylmethyl ester CAS No.:102-16-9 Molecular Structure: Molecular Structure of 102-16-9 (Benzeneacetic acid,phenylmethyl ester) Formula: C15H14O2 Molecular Weight : 226.27 Synonyms: BENZYL PHENYLACETATE;Phenylacetic Acid Benzyl Ester;benzyl 2-phenylacetate; EINECS: 203-008-4 Density: 1.147

-

7-BROMO-1-NITRONAPHTHALENE CAS 102153-49-1

Name: Naphthalene, 7-bromo-1-nitro- CAS No.:102153-49-1 Molecular Structure: Molecular Structure of 102153-49-1 (Naphthalene, 7-bromo-1-nitro-) Formula: C10H6BrNO2 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport

- ## 2-(2,2-Dimethyl-1,3-dioxan-5-yl)ethanol CAS 102147-75-1

  Name: 1,3-Dioxane-5-ethanol, 2,2-dimethyl- CAS No.:102147-75-1 Molecular Structure: Molecular Structure of 102147-75-1 (1,3-Dioxane-5-ethanol, 2,2-dimethyl-) Formula: C8H16O3 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes:

- ## SODIUM METASILICATE PENTAHYDRATE CAS 10213-79-3

  Name: Sodium metasilicate pentahydrate CAS No.:10213-79-3 Molecular Structure: Molecular Structure of 10213-79-3 (Sodium metasilicate pentahydrate) Formula: H10Na2O8Si Molecular Weight : 212.14 Synonyms: Disodiumsilicate pentahydrate (Na2SiO3.5H2O);Dorimetakeiso 5aq;Dry Metakeiso 5aq;Metaesu5;Metso granular;Sodium

- ## Sodium tungstate dihydrate CAS 10213-10-2

  Name: Sodium tungstate dihydrate CAS No.:10213-10-2 Molecular Structure: Molecular Structure of 10213-10-2 (Sodium tungstate dihydrate) Formula: H2O.Na.1/2O4W Molecular Weight : 164.62 Synonyms: Tungsticacid (H2WO4), disodium salt, dihydrate (8CI);Disodium tungstate (Na2WO4)dihydrate;Sodium tungstate

- ## Piperazine, 1-(cyclopropylcarbonyl)-, Monohydrochloride CAS 1021298-67-8

  Name: Piperazine, 1-(cyclopropylcarbonyl)-, Monohydrochloride CAS No.:1021298-67-8 Molecular Structure: Molecular Structure of 1021298-67-8 (Piperazine, 1-(cyclopropylcarbonyl)-, Monohydrochloride) Formula: Molecular Weight : 190.6705 Synonyms: Piperazine, 1-(cyclopropylcarbonyl)-,

- ## (3S)-3-(tert-Butoxycarbonyl)amino-1-chloro-4-phenyl-2-butanone CAS 102123-74-0

  Name: (3S)-3-(tert-Butoxycarbonyl)amino-1-chloro-4-phenyl-2-butanone CAS No.:102123-74-0 Molecular Structure: Molecular Structure of 102123-74-0 ((3S)-3-(tert-Butoxycarbonyl)amino-1-chloro-4-phenyl-2-butanone) Formula: C15H20ClNO3 Molecular Weight : 297.78 Synonyms: Carbamicacid,

- ## 4-Amino-1-[(2R,3R,4R,5R)-3-fluoro-4-hydroxy-5-(hydroxymethyl)oxolan-2-yl]pyrimidin-2-one CAS 10212-20-1

  Name: Cytidine,2'-deoxy-2'-fluoro- CAS No.:10212-20-1 Molecular Structure: Molecular Structure of 10212-20-1 (Cytidine,2'-deoxy-2'-fluoro-) Formula: C9H12FN3O4 Molecular Weight : 245.21 Synonyms:

- ## AZOCASEIN CAS 102110-74-7

  Name: Caseins, azo CAS No.:102110-74-7 Molecular Structure: Molecular Structure of 102110-74-7 (Caseins, azo) Formula: Molecular Weight : 0 Synonyms: Azo caseins EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

- ## Tin ores, concs., leached CAS 102110-65-6

  Name: Tin ores, concs.,leached CAS No.:102110-65-6 Molecular Structure: Molecular Structure of 102110-65-6 (Tin ores, concs.,leached) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

- ## 2-Acetylanthracene CAS 10210-32-9

  Name: Ethanone,1-(2-anthracenyl)- CAS No.:10210-32-9 Molecular Structure: Molecular Structure of 10210-32-9 (Ethanone,1-(2-anthracenyl)-) Formula: C16H12O Molecular Weight : 220.27 Synonyms: Ketone,2-anthryl methyl (7CI,8CI);2-Anthryl methyl ketone; EINECS: 227-730-4 Density: 1.164 g/cm3 Melting Point: Boiling Point:

- ## α-muurolene CAS 10208-80-7
  Name: Naphthalene,1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-(1-methylethyl)-,(57264155,1S,4aS,8aR)- CAS No.:10208-80-7 Molecular Structure: Molecular Structure of 10208-80-7 (Naphthalene,1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-(1-methylethyl)-,(57264156,1S,4aS,8aR)-) Formula: C15H24 Molecular Weight : 0 Synonyms:

- ## N,N'-Diphenylurea CAS 102-07-8
  Name: Urea, N,N'-diphenyl- CAS No.:102-07-8 Molecular Structure: Molecular Structure of 102-07-8 (Urea, N,N'-diphenyl-) Formula: C13H12N2O Molecular Weight : 212.27 Synonyms: Carbanilide(7CI,8CI);1,3-Diphenylcarbamide;AD 30;DPU;N,N'-Diphenylurea;N-Phenyl-N'-phenylurea;NSC 227401;NSC

- ## (5-METHYLPYRIDIN-3-YL)METHANOL CAS 102074-19-1
  Name: 3-Pyridinemethanol,5-methyl- CAS No.:102074-19-1 Molecular Structure: Molecular Structure of 102074-19-1 (3-Pyridinemethanol,5-methyl-) Formula: C7H9NO Molecular Weight : 123.15 Synonyms: (5-Methyl-3-pyridyl)methan-1-ol;3-Methylpyridine-5-methanol; EINECS: Density: 1.092 g/cm3 Melting Point: Boiling Point:

- ## 99% CAS 1020665-14-8
  Name: Neuburg siliceous earth CAS No.:1020665-14-8 Molecular Structure: Molecular Structure of 1020665-14-8 (Neuburg siliceous earth) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

- ## DIBEMETHINE CAS 102-05-6
  Name: Benzenemethanamine,N-methyl-N-(phenylmethyl)- CAS No.:102-05-6 Molecular Structure: Molecular Structure of 102-05-6 (Benzenemethanamine,N-methyl-N-(phenylmethyl)-) Formula: C15H17N Molecular Weight : 211.30 Synonyms: Dibenzylamine,N-methyl-

- ## BIS(1-METHOXY-2-PROPYL)MALEATE CAS 102054-10-4
  Name: 2-Butenedioic acid(2Z)-, 1,4-bis(2-methoxy-1-methylethyl) ester CAS No.:102054-10-4 Molecular Structure: Molecular Structure of 102054-10-4 (2-Butenedioic acid(2Z)-, 1,4-bis(2-methoxy-1-methylethyl) ester) Formula: C12H20 O6 Molecular Weight : 260.28 Synonyms: 2-Butenedioicacid (2Z)-,

- ## 1,3-Diphenylacetone CAS 102-04-5
  Name: 1,3-Diphenylacetone CAS No.:102-04-5 Molecular Structure: Molecular Structure of 102-04-5 (1,3-Diphenylacetone) Formula: C15H14O Molecular Weight : 210.28 Synonyms: 1,3-Diphenyl-2-propanone;2-Propanone,1,3-diphenyl-;1,3-Diphenylpropanone;Benzyl ketone;Dibenzyl ketone;NSC220312;NSC 407392;NSC

- ## Diethyl isobutylmalonate CAS 10203-58-4
  Name: Propanedioic acid,2-(2-methylpropyl)-, 1,3-diethyl ester CAS No.:10203-58-4 Molecular Structure: Molecular Structure of 10203-58-4 (Propanedioic acid,2-(2-methylpropyl)-, 1,3-diethyl ester) Formula: C11H20O4 Molecular Weight : 216.28 Synonyms: Malonicacid, isobutyl-, diethyl ester (6CI,7CI,8CI);Propanedioic

- ## 3,5-Dichlorobenzaldehyde CAS 10203-08-4
  Name: Benzaldehyde, 3,5-dichloro- CAS No.:10203-08-4 Molecular Structure: Molecular Structure of 10203-08-4 (Benzaldehyde, 3,5-dichloro-) Formula: C7H4Cl2O Molecular Weight : 175.01 Synonyms: NSC 109095; EINECS: 233-499-0 Density: 1.4 g/cm3 Melting Point: Boiling Point: 241.1 °C at 760 mmHg Flash Point: 98.3 °C

- ## C53031900 CAS 1020251-53-9
  Name: Benzamide, N-[(3,4-dimethoxyphenyl)methyl]-2-[(2-hydroxy-1-methylethyl)amino]-5-nitro- CAS No.:1020251-53-9 Molecular Structure: Molecular Structure of 1020251-53-9 (Benzamide, N-[(3,4-dimethoxyphenyl)methyl]-2-[(2-hydroxy-1-methylethyl)amino]-5-nitro-) Formula: C19H23N3O6 Molecular Weight : 389.4057 Synonyms:

- ## MeayaMycin B CAS 1020210-12-1
  0.99

- ## 99% CAS 10200-71-2
  0.99

- ## 1,3-Thiazole-2-carbaldehyde CAS 10200-59-6
  Name: 2-Thiazolecarboxaldehyde CAS No.:10200-59-6 Molecular Structure: Molecular Structure of 10200-59-6 (2-Thiazolecarboxaldehyde) Formula: C4H3NOS Molecular Weight : 113.14 Synonyms: thiazole-2-carbaldehyde;Thiazole-2-carboxaldehyde;1,3-Thiazole-2-carboxaldehyde;2-Formylthiazole; EINECS: Density: 1.288 g/cm3 Melting

- ## 99% CAS 10200-48-3
  Name: 1,2-Oxathiolan-4-ol, 2,2-dioxide CAS No.:10200-48-3 Molecular Structure: Molecular Structure of 10200-48-3 (1,2-Oxathiolan-4-ol, 2,2-dioxide) Formula: C3H6O4S Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport

- ## 4'-PHOSPHOPANTETHEINE CAS 1019842-99-9
  Name: 4'-Phosphopantetheine CAS No.:1019842-99-9 Molecular Structure: Molecular Structure of 1019842-99-9 (4'-Phosphopantetheine) Formula: C11H23N2O7PS Molecular Weight : 358.35 Synonyms: PPT;4'-PHOSPHOPANTETHEINE;2,8-bis(diphenylphosphoryl)dibenzo[b,d]thiophene;PPT , 2,8-bis(diphenylphosphoryl)dibenzo[b,d]thiophene

- ## N-Benzyl-N-methylethanolamine CAS 101-98-4
  Name: Ethanol, 2-[methyl(phenylmethyl)amino]- CAS No.:101-98-4 Molecular Structure: Molecular Structure of 101-98-4 (Ethanol, 2-[methyl(phenylmethyl)amino]-) Formula: C10H15NO Molecular Weight : 165.23 Synonyms: Ethanol,2-(benzylmethylamino)-

- ## Chlorodi(p-tolyl)phosphine, 95% CAS 1019-71-2
  Name: Chloro di(p-tolyl)phosphine CAS No.:1019-71-2 Molecular Structure: Molecular Structure of 1019-71-2 (Chloro di(p-tolyl)phosphine) Formula: C14H14ClP Molecular Weight : 248.69 Synonyms: Di(p-tolyl)chlorophosphine; EINECS: Density: 1.1 g/mL at 25 °C Melting Point: Boiling Point: 346.9 °C at 760 mmHg Flash Point:

- ## tetrahydroxysilane CAS 10193-36-9
  Name: Silicic acid (H4O4Si) CAS No.:10193-36-9 Molecular Structure: Molecular Structure of 10193-36-9 (Silicic acid (H4O4Si)) Formula: H4O4Si Molecular Weight : 96.11486 Synonyms: Monosilicicacid;Orthosilicic acid;Orthosilicic acid (H4SiO4);Silicic acid (Si(OH)4);Silicon hydroxide (Si(OH)4);Silicon

- ## 4-[4-[[[[4-Chloro-3-(trifluoromethyl)phenyl]amino]carbonyl]amino]-3-fluorophenoxy]-N-methyl-2-pyridinecarboxamide hydrate CAS 1019206-88-2
  Name: Regorafenib hydrate CAS No.:1019206-88-2 Molecular Structure: Molecular Structure of 1019206-88-2 (Regorafenib hydrate) Formula: C21H15ClF4N4O3 Molecular Weight : Synonyms: Regorafenib monohydrate EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes:

N-Cyanoimido-S,S-dimethyl-dithiocarbonate CAS 10191-60-3
Name: Carbonimidodithioic acid,N-cyano-,dimethyl ester CAS No.:10191-60-3 Molecular Structure: Molecular Structure of 10191-60-3 (Carbonimidodithioic acid,N-cyano-,dimethyl ester) Formula: C4H6N2S2 Molecular Weight : 146.22 Synonyms: Carbonimidodithioicacid, cyano-, dimethyl ester (9CI);Imidocarbonic acid,

- ## LEAD MOLYBDATE CAS 10190-55-3
  Name: Lead molybdate CAS No.:10190-55-3 Molecular Structure: Molecular Structure of 10190-55-3 (Lead molybdate) Formula: MoO4Pb Molecular Weight : 367.14 Synonyms: LEAD MOLYBDATE;LEAD MOLYBDENUM OXIDE;LEAD(II) MOLYBDATE;leadmolybdenumoxide(pbmoo4);Lead (II) molybdate (99.95+% Pb) (metals basis);LEAD(II) MOLYBDATE,

- ## Distillates (coal tar), benzole fraction, BTX-rich CAS 101896-26-8
  Name: Distillates (coal tar),benzole fraction, BTX-rich CAS No.:101896-26-8 Molecular Structure: Molecular Structure of 101896-26-8 (Distillates (coal tar),benzole fraction, BTX-rich) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard

- ## PYRROLNITRIN CAS 1018-71-9
  Name: 1H-Pyrrole,3-chloro-4-(3-chloro-2-nitrophenyl)- CAS No.:1018-71-9 Molecular Structure: Molecular Structure of 1018-71-9 (1H-Pyrrole,3-chloro-4-(3-chloro-2-nitrophenyl)-) Formula: C10H6 Cl2 N2 O2 Molecular Weight : 257.08 Synonyms: Pyrrole,3-chloro-4-(3-chloro-2-nitrophenyl)-

- ## Diclazuril CAS 101831-37-2
  Name: Diclazuril CAS No.:101831-37-2 Molecular Structure: Molecular Structure of 101831-37-2 (Diclazuril) Formula: C17H9Cl3N4O2 Molecular Weight : 407.64 Synonyms: P 64433;2,6-Dichloro-.alpha.-(4-chlorophenyl)-4-(4,5-dihydro-3,5-dioxo-1,2,4-triazin-2(3H)- yl)benzeneacetonitrile;Benzeneacetonitrile,

- ## 4,4'-Oxydianiline CAS 101-80-4
  Name: 4,4'-Oxydianiline CAS No.:101-80-4 Molecular Structure: Molecular Structure of 101-80-4 (4,4'-Oxydianiline) Formula: C12H12N2O Molecular Weight : 200.23648 [g/mol] Synonyms: 4,4'-Diaminobiphenyl oxide;4,4'-Diaminodiphenyl ether;4,4'-Diaminodiphenyl

- ## 4,5-Difluoro-m-toluic acid, 5-Carboxy-2,3-difluorotoluene CAS 1017778-60-7
  Name: 4,5-Difluoro-m-toluic acid, 5-Carboxy-2,3-difluorotoluene CAS No.:1017778-60-7 Molecular Structure: Molecular Structure of 1017778-60-7 (4,5-Difluoro-m-toluic acid, 5-Carboxy-2,3-difluorotoluene) Formula: Molecular Weight : 172.1288464 Synonyms: 3,4-Difluoro-5-methylbenzoic acid;4,5-Difluoro-m-toluic acid,

- ## 4,4'-DIMETHOXYDIPHENYLAMINE CAS 101-70-2
  Name: Benzenamine,4-methoxy-N-(4-methoxyphenyl)- CAS No.:101-70-2 Molecular Structure: Molecular Structure of 101-70-2 (Benzenamine,4-methoxy-N-(4-methoxyphenyl)-) Formula: C14H15NO2 Molecular Weight : 229.30 Synonyms: Diphenylamine,4,4'-dimethoxy-

- ## (1R,2R,3aS,9aS)-2,3,3a,4,9,9a-Hexahydro-1-[(3S)-3-hydroxyoctyl]-1H-benz[f]indene-2,5-diol CAS 101692-02-8
  0.99

- ## ACID RED 97 (C.I. 22890) CAS 10169-02-5
  Name: [1,1'-Biphenyl]-2,2'-disulfonicacid, 4,4'-bis[2-(2-hydroxy-1-naphthalenyl)diazenyl]-, sodium salt (1:2) CAS No.:10169-02-5 Molecular Structure: Molecular Structure of 10169-02-5 ([1,1'-Biphenyl]-2,2'-disulfonicacid, 4,4'-bis[2-(2-hydroxy-1-naphthalenyl)diazenyl]-, sodium salt (1:2)) Formula: C32H20N4O8S2.2Na

- ## 4,4'-Diphenylmethane diisocyanate CAS 101-68-8
  Name: Benzene,1,1'-methylenebis[4-isocyanato- CAS No.:101-68-8 Molecular Structure: Molecular Structure of 101-68-8 (Benzene,1,1'-methylenebis[4-isocyanato-) Formula: C15H10N2O2 Molecular Weight : 250.27 Synonyms: 4,4'-Methylenediphenyl diisocyanate;4,4'-Methylenediphenylene

99% CAS 101685-29-4
Name: 1-Propene-1,1,3-tricarbonitrile, 2-phenyl- CAS No.:101685-29-4 Molecular Structure: Molecular Structure of 101685-29-4 (1-Propene-1,1,3-tricarbonitrile, 2-phenyl-) Formula: C12H7N3 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols:

5-METHOXY-PYRIDIN-2-YLAMINE CAS 10167-97-2
Name: 2-Pyridinamine,5-methoxy- CAS No.:10167-97-2 Molecular Structure: Molecular Structure of 10167-97-2 (2-Pyridinamine,5-methoxy-) Formula: C6H8N2O Molecular Weight : 124.14052 Synonyms: Pyridine,2-amino-5-methoxy- (7CI,8CI);5-Methoxy-2-aminopyridine;5-Methoxypyridin-2-amine; EINECS: Density: 1.139 g/cm3 Melting

5-METHYL-4-OXO-3,4-DIHYDRO-THIENO[2,3-D]PYRIMIDINE-6-CARBOXYLIC ACID CAS 101667-97-4

Name: 7-Methyl-5-oxo-9-thia-2,4-diazabicyclo[4.3.0]nona-2,7,10-triene-8-carboxylate CAS No.:101667-97-4 Molecular Structure: Molecular Structure of 101667-97-4 (7-Methyl-5-oxo-9-thia-2,4-diazabicyclo[4.3.0]nona-2,7,10-triene-8-carboxylate) Formula: C8H6N2O3S Molecular Weight : 210.21 Synonyms:

Silicic acid, magnesium sodium salt CAS 101659-01-2

Name: Silicic acid, magnesiumsodium salt CAS No.:101659-01-2 Molecular Structure: Molecular Structure of 101659-01-2 (Silicic acid, magnesiumsodium salt) Formula: Molecular Weight : 0 Synonyms: LaponiteXL; Magnesium sodium silicate; Sodium magnesium silicate EINECS: Density: Melting Point: Boiling Point: Flash Point:

4,5-DIMETHOXY-2-NITROBENZYL ALCOHOL CAS 1016-58-6

Name: Benzenemethanol,4,5-dimethoxy-2-nitro- CAS No.:1016-58-6 Molecular Structure: Molecular Structure of 1016-58-6 (Benzenemethanol,4,5-dimethoxy-2-nitro-) Formula: C9H11NO5 Molecular Weight : 213.1873 Synonyms: Veratrylalcohol, 6-nitro- (7CI,8CI);2-Nitro-4,5-dimethoxybenzyl alcohol;4,5-Dimethoxy-2-nitrobenzyl

4,4'-DINITRODIPHENYL ETHER CAS 101-63-3

Name: Benzene,1,1'-oxybis[4-nitro- CAS No.:101-63-3 Molecular Structure: Molecular Structure of 101-63-3 (Benzene,1,1'-oxybis[4-nitro-) Formula: C12H8N2O5 Molecular Weight : 260.22 Synonyms: Ether,bis(p-nitrophenyl) (6CI,7CI,8CI);4,4'-Dinitrodiphenyl oxide;Bis(4-nitrophenyl) ether;Bis(p-nitrophenyl)

99% CAS 101632-25-9

0.99

TALIPEXOLE CAS 101626-70-4

Name: 4H-Thiazolo[4,5-d]azepin-2-amine,5,6,7,8-tetrahydro-6-(2-propen-1-yl)- CAS No.:101626-70-4 Molecular Structure: Molecular Structure of 101626-70-4 (4H-Thiazolo[4,5-d]azepin-2-amine,5,6,7,8-tetrahydro-6-(2-propen-1-yl)-) Formula: C10H15 N3 S Molecular Weight : 209.31 Synonyms:

1,3-Dibromo-2-methyl-5-nitrobenzene CAS 101581-06-0

Name: Benzene, 1,3-dibromo-2-methyl-5-nitro- CAS No.:101581-06-0 Molecular Structure: Molecular Structure of 101581-06-0 (Benzene, 1,3-dibromo-2-methyl-5-nitro-) Formula: C7H5Br2NO2 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk

Dodecyl 4-methylbenzenesulfonate CAS 10157-76-3

Name: Benzenesulfonic acid,4-methyl-, dodecyl ester CAS No.:10157-76-3 Molecular Structure: Molecular Structure of 10157-76-3 (Benzenesulfonic acid,4-methyl-, dodecyl ester) Formula: C19H32O3S Molecular Weight : 340.52 Synonyms: Dodecylalcohol, p-toluenesulfonate (6CI);p-Toluenesulfonic acid, dodecyl

Fmoc-D-proline CAS 101555-62-8

Name: 1,2-Pyrrolidinedicarboxylicacid, 1-(9H-fluoren-9-ylmethyl) ester,(57264109,2R)- CAS No.:101555-62-8 Molecular Structure: Molecular Structure of 101555-62-8 (1,2-Pyrrolidinedicarboxylicacid, 1-(9H-fluoren-9-ylmethyl) ester,(57264110,2R)-) Formula: C20H19NO4 Molecular Weight : 337.37 Synonyms:

BOC-D-BETA-HOMOPROLINE CAS 101555-60-6

Name: 2-Pyrrolidineaceticacid, 1-[(1,1-dimethylethoxy)carbonyl]-,(57264104,2R)- CAS No.:101555-60-6 Molecular Structure: Molecular Structure of 101555-60-6 (2-Pyrrolidineaceticacid, 1-[(1,1-dimethylethoxy)carbonyl]-,(57264105,2R)-) Formula: C11H19 N O4 Molecular Weight : 229.27 Synonyms: 2-Pyrrolidineaceticacid,

3-formyl-5-hydroxybenzonitrile CAS 1015414-78-4

0.99

DIPHENYLPHOSPHINOTHIOYL CHLORIDE CAS 1015-37-8

Name: Phosphinothioicchloride, P,P-diphenyl- CAS No.:1015-37-8 Molecular Structure: Molecular Structure of 1015-37-8 (Phosphinothioicchloride, P,P-diphenyl-) Formula: C12H10 Cl P S Molecular Weight : 252.7 Synonyms: Phosphinothioicchloride, diphenyl- (6CI,7CI,8CI,9CI); Chlorodiphenylphosphine

3-Chloro-2,4,5-trifluorobenzoic acid CAS 101513-77-3

Name: Benzoic acid,3-chloro-2,4,5-trifluoro- CAS No.:101513-77-3 Molecular Structure: Molecular Structure of 101513-77-3 (Benzoic acid,3-chloro-2,4,5-trifluoro-) Formula: C7H2ClF3O2 Molecular Weight : 210.5378 Synonyms: 2,4,5-Trifluoro-3-chlorobenzoic acid; EINECS: Density: 1.663 g/cm3 Melting Point: Boiling Point:

1,4-Bis-(isocyanatomethyl)-benzene CAS 1014-98-8

Name: Benzene,1,4-bis(isocyanatomethyl)- CAS No.:1014-98-8 Molecular Structure: Molecular Structure of 1014-98-8 (Benzene,1,4-bis(isocyanatomethyl)-) Formula: C10H8 N2 O2 Molecular Weight : 186.18 Synonyms: Isocyanicacid, p-phenylenedimethylene ester (7CI,8CI); 1,4-Bis(isocyanatomethyl)benzene;1,4-Xylylene

1-Propanesulfonyl chloride CAS 10147-36-1

Name: 1-Propanesulfonyl chloride CAS No.:10147-36-1 Molecular Structure: Molecular Structure of 10147-36-1 (1-Propanesulfonyl chloride) Formula: C3H7ClO2S Molecular Weight : 142.61 Synonyms: Propanesulfonylchloride (6CI);Propylsulfonyl chloride;n-Propanesulfonyl chloride;n-Propylsulphonyl chloride; EINECS: 233-414-7

FENURON CAS 101-42-8

Name: 1,1-Dimethyl-3-phenylurea CAS No.:101-42-8 Molecular Structure: Molecular Structure of 101-42-8 (1,1-Dimethyl-3-phenylurea) Formula: C9H12 N2 O Molecular Weight : 164.23 Synonyms: Urea,1,1-dimethyl-3-phenyl- (8CI); 1,1-Dimethyl-3-phenylurea;1-Phenyl-3,3-dimethylurea; 3-Phenyl-1,1-dimethylurea; Amicure UR; Dibar;

Alizarine Sky Blue 3FG CAS 10142-59-3

Name: Benzoic acid,5-[[[3-[(4-amino-9,10-dihydro-9,10-dioxo-3-sulfo-1-anthracenyl)amino]-4-methoxyphenyl]methyl]sulfonyl]-2-hydroxy-,sodium salt (1:2) CAS No.:10142-59-3 Molecular Structure: Molecular Structure of 10142-59-3 (Benzoic

CHROMIUM POTASSIUM SULFATE CAS 10141-00-1

Name: Sulfuric acid,chromium(3+) potassium salt (2:1:1) CAS No.:10141-00-1 Molecular Structure: Molecular Structure of 10141-00-1 (Sulfuric acid,chromium(3+) potassium salt (2:1:1)) Formula: Cr. 2 H2 O4 S . K Molecular Weight : 283.22 Synonyms: Chromealum; Chrome potash alum; Chromic potassium sulfate; Chromium

SODIUM PERBORATE TETRAHYDRATE CAS 10140-63-3

Name: Boric acid (H3BO3),monosodium salt, tetrahydrate (8CI,9CI) CAS No.:10140-63-3 Molecular Structure: Molecular Structure of 10140-63-3 (Boric acid (H3BO3),monosodium salt, tetrahydrate (8CI,9CI)) Formula: BH3 O3 . 4 H2 O . Na Molecular Weight : 153.86 Synonyms: SODIUM PEROXOBORATE TETRAHYDRATE;SODIUM

Fatty acids, palm kernel-oil CAS 101403-98-9

Name: Fatty acids, palm kernel-oil CAS No.:101403-98-9 Molecular Structure: Molecular Structure of 101403-98-9 (Fatty acids, palm kernel-oil) Formula: Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information:

alpha-Methylcinnamaldehyde CAS 101-39-3

Name: 2-Propenal,2-methyl-3-phenyl- CAS No.:101-39-3 Molecular Structure: Molecular Structure of 101-39-3 (2-Propenal,2-methyl-3-phenyl-) Formula: C10H10O Molecular Weight : 146.19 Synonyms: Cinnamaldehyde,a-methyl-

2-Chloro-1,3-dimethylimidazolidinium hexafluorophosphate CAS 101385-69-7

Name: 2-Chloro-1,3-dimethylimidazolidinium hexafluorophosphate CAS No.:101385-69-7 Molecular Structure: Molecular Structure of 101385-69-7 (2-Chloro-1,3-dimethylimidazolidinium hexafluorophosphate) Formula: C5H10ClF6N2P Molecular Weight : 278.56 Synonyms: 2-CHLORO-1,3-DIMETHYLIMIDAZOLINIUM HEXAFLUOROPHOSPHATE;CIP

Ammonium iron(III) sulfate CAS 10138-04-2

Name: Sulfuric acid, ammoniumiron(3+) salt (2:1:1) CAS No.:10138-04-2 Molecular Structure: Molecular Structure of 10138-04-2 (Sulfuric acid, ammoniumiron(3+) salt (2:1:1)) Formula: Fe.NH4.(SO4)2 Molecular Weight : 266.01 Synonyms: Ammoniumiron sulfate (NH4Fe(SO4)2) (6CI);Ammonium ferric alum;Ammonium

(E,Z,Z)-4,6,10-Hexadecatrien-1-ol CAS 101373-00-6

Name: (E,Z,Z)-4,6,10-Hexadecatrien-1-ol CAS No.:101373-00-6 Molecular Structure: Molecular Structure of 101373-00-6 ((E,Z,Z)-4,6,10-Hexadecatrien-1-ol) Formula: Molecular Weight : 236.39292 Synonyms: (E,Z,Z)-4,6,10-Hexadecatrien-1-ol;4E6Z10Z-16OH;(4E,6Z,10Z)-hexadeca-4,6,10-trien-1-ol EINECS: Density: Melting Point:

2,6-bis(3-(9H-carbazol-9-yl)phenyl)pyridine CAS 1013405-24-7

Name: 9H-Carbazole, 9,9'-(2,6-pyridinediyldi-3,1-phenylene)bis- CAS No.:1013405-24-7 Molecular Structure: Molecular Structure of 1013405-24-7 (9H-Carbazole, 9,9'-(2,6-pyridinediyldi-3,1-phenylene)bis-) Formula: C41H27N3 Molecular Weight : 561.67 Synonyms: 2,6-Bis[3-(carbazol-9-yl)phenyl]pyridine; EINECS: Density: 1.21

BENZO[D]OXAZOL-2-YLMETHANAMINE CAS 101333-98-6

Name: 2-Aminomethylbenzooxazole CAS No.:101333-98-6 Molecular Structure: Molecular Structure of 101333-98-6 (2-Aminomethylbenzooxazole) Formula: C8H8N2O Molecular Weight : 148.16 Synonyms: 2-Benzoxazolemethanamine;1,3-Benzoxazol-2-ylmethylamine; EINECS: Density: 1.239 g/cm3 Melting Point: Boiling Point: 246.651 °C at

3-BROMOPROPIONIC ACID N-HYDROXYSUCCINIMIDE CAS 101314-84-5

Name: 3-BROMOPROPIONIC ACID N-HYDROXYSUCCINIMIDE CAS No.:101314-84-5 Molecular Structure: Molecular Structure of 101314-84-5 (3-BROMOPROPIONIC ACID N-HYDROXYSUCCINIMIDE) Formula: C7H12BrNO4 Molecular Weight : 254.07848 Synonyms: 3-BROMOPROPIONIC ACID

Trans-1-benzyl-4-(4-chlorophenyl)pyrrolidine-3-carboxylic acid-HCl CAS 1013117-42-4

Name: Trans-1-benzyl-4-(4-chlorophenyl)pyrrolidine-3-carboxylic acid-HCl CAS No.:1013117-42-4 Molecular Structure: Molecular Structure of 1013117-42-4 (Trans-1-benzyl-4-(4-chlorophenyl)pyrrolidine-3-carboxylic acid-HCl) Formula: Molecular Weight : 352.25496 Synonyms: EINECS: Density: Melting Point: Boiling Point:

99% CAS 101289-18-3

0.99

ACID ORANGE 74 CAS 10127-27-2

Name: Chromate(1-),[3-[2-[4,5-dihydro-3-methyl-5-(oxo-kO)-1-phenyl-1H-pyrazol-4-yl]diazenyl-kN1]-2-(hydroxy-kO)-5-nitrobenzenesulfonato(3-)]hydroxy-, sodium (1:1),(57264078,T-4)- CAS No.:10127-27-2 Molecular Structure: Molecular Structure of 10127-27-2

1,4-Di-tert-butylbenzene CAS 1012-72-2

Name: 1,4-Di-tert-butylbenzene CAS No.:1012-72-2 Molecular Structure: Molecular Structure of 1012-72-2 (1,4-Di-tert-butylbenzene) Formula: C14H22 Molecular Weight : 190.36 Synonyms: Benzene,p-di-tert-butyl-(7CI,8CI);1,4-Bis(1,1-dimethylethyl)benzene;1,4-Bis(tert-butyl)benzene;NSC

4,4'-[(3,3'-dimethoxy[1,1'-biphenyl]-4,4'-diyl)bis(azo)]bis[3-hydroxy-N-phenylnaphthalene-2-carboxamide] CAS 10127-03-4

Name: 2-Naphthalenecarboxamide,4,4'-[(3,3'-dimethoxy[1,1'-biphenyl]-4,4'-diyl)bis(2,1-diazenediyl)]bis[3-hydroxy-N-phenyl- CAS No.:10127-03-4 Molecular Structure: Molecular Structure of 10127-03-4 (2-Naphthalenecarboxamide,4,4'-[(3,3'-dimethoxy[1,1'-biphenyl]-4,4'-diyl)bis(2,1-diazenediyl)]bis[3-hydroxy-N-phenyl-)

Basic Orange 14 CAS 10127-02-3

Name: Euchrysine 3RX CAS No.:10127-02-3 Molecular Structure: Molecular Structure of 10127-02-3 (Euchrysine 3RX) Formula: C17H19N3.xCl2Zn.ClH Molecular Weight : 739.91400 Synonyms: 3,6-Acridinediamine,N,N,N',N'-tetramethyl-,monohydrochloride,compd. with zinc chloride (ZnCl2);Basic Orange 14;Acridine Orange, hemi(zinc

Mestinon CAS 101-26-8

Name: Mestinon CAS No.:101-26-8 Molecular Structure: Molecular Structure of 101-26-8 (Mestinon) Formula: C9H13BrN2O2 Molecular Weight : 261.15 Synonyms: Pyridinium, 3-hydroxy-1-methyl-, bromide, dimethylcarbamate(8CI);Carbamic acid, dimethyl-, ester with 3-hydroxy-1-methylpyridiniumbromide

4-heptadecylidene-3-hexadecyloxetan-2-one CAS 10126-68-8

Name: 2-Oxetanone,4-heptadecylidene-3-hexadecyl- CAS No.:10126-68-8 Molecular Structure: Molecular Structure of 10126-68-8 (2-Oxetanone,4-heptadecylidene-3-hexadecyl-) Formula: C36H68O2 Molecular Weight : 532.92392 Synonyms: 3-Eicosenoicacid, 2-hexadecyl-3-hydroxy-, b-lactone

Copper(II) chloride dihydrate CAS 10125-13-0

Name: Copper(II) chloride dihydrate CAS No.:10125-13-0 Molecular Structure: Molecular Structure of 10125-13-0 (Copper(II) chloride dihydrate) Formula: CuCl2.2(H2O) Molecular Weight : 170.48 Synonyms: Copperchloride (CuCl2.2H2O);Copper dichloride dihydrate;Copper(II) chloride dihydrate;Cupric chloride dihydrate;

99% CAS 101250-97-9

Name: Guanidine, N,N'-dimethyl-N''-nitro- CAS No.:101250-97-9 Molecular Structure: Molecular Structure of 101250-97-9 (Guanidine, N,N'-dimethyl-N''-nitro-) Formula: C3H8N4O2 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes:

COPPERSULFATE CAS 10124-44-4

Name: Sulfuric acid, coppersalt (8CI,9CI) CAS No.:10124-44-4 Molecular Structure: Molecular Structure of 10124-44-4 (Sulfuric acid, coppersalt (8CI,9CI)) Formula: CuO4S Molecular Weight : 159.61 Synonyms: Sulfuric acid, copper salt; EINECS: 231-847-6;215-568-7 Density: Melting Point: Boiling Point: 330°C at 760 mmHg

TRI-AMMONIUM ORTHOPHOSPHATE CAS 10124-31-9

Name: Phosphoric acid,ammonium salt (1:?) CAS No.:10124-31-9 Molecular Structure: Molecular Structure of 10124-31-9 (Phosphoric acid,ammonium salt (1:?)) Formula: H3N . x H3O4P Molecular Weight : 149.09 Synonyms: Phosphoricacid, ammonium salt (8CI,9CI);Ammonium acid phosphate;Ammoniumorthophosphate;Fyrex;Nonnen SMC

(2R,4S)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpentanoic acid CAS 1012341-50-2

Name: (2R,4S)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpentanoic acid CAS No.:1012341-50-2 Molecular Structure: Molecular Structure of 1012341-50-2 ((2R,4S)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpentanoic acid) Formula: Molecular Weight : 383.48066 Synonyms:

(R,E)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpent-2-enoic acid CAS 1012341-48-8

Name: (R,E)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpent-2-enoic acid CAS No.:1012341-48-8 Molecular Structure: Molecular Structure of 1012341-48-8 ((R,E)-5-([1,1'-biphenyl]-4-yl)-4-((tert-butoxycarbonyl)aMino)-2-Methylpent-2-enoic acid) Formula: Molecular Weight : 381.46478 Synonyms:

4-(TRIMETHYLSILOXY)BENZALDEHYDE CAS 1012-12-0

Name: Benzaldehyde,4-[(trimethylsilyl)oxy]- CAS No.:1012-12-0 Molecular Structure: Molecular Structure of 1012-12-0 (Benzaldehyde,4-[(trimethylsilyl)oxy]-) Formula: C10H14O2Si Molecular Weight : 194.3 Synonyms: Benzaldehyde,p-(trimethylsiloxy)-

DL-Homophenylalanine CAS 1012-05-1

Name: Benzenebutanoic acid, a-amino- CAS No.:1012-05-1 Molecular Structure: Molecular Structure of 1012-05-1 (Benzenebutanoic acid, a-amino-) Formula: C10H13NO2 Molecular Weight : 179.22 Synonyms: Butyric acid,2-amino-4-phenyl-, DL- (8CI);2-Amino-4-phenylbutanoic

SODIUM PHENOTHIAZINE-10-YL-PROPYLSULFONATE CAS 101199-38-6

Name: 10H-Phenothiazine-10-propanesulfonicacid, sodium salt (1:1) CAS No.:101199-38-6 Molecular Structure: Molecular Structure of 101199-38-6 (10H-Phenothiazine-10-propanesulfonicacid, sodium salt (1:1)) Formula: C15H14NNaO3S2 Molecular Weight : 343.03 Synonyms: 10H-Phenothiazine-10-propanesulfonicacid, sodium salt

Methyl indole-5-carboxylate CAS 1011-65-0

Name: Methyl indole-5-carboxylate CAS No.:1011-65-0 Molecular Structure: Molecular Structure of 1011-65-0 (Methyl indole-5-carboxylate) Formula: C10H9NO2 Molecular Weight : 175.184 Synonyms: 5-Indolecarboxylic Acid Acid Methyl Ester;Indole-5-carboxylic acid methyl ester;1H-Indole-5-carboxylic acid methyl ester;methyl

(1R,2S)-rel-2-(Aminomethyl)-N,N-diethyl-1-phenylcyclopropanecarboxamide hydrochloride CAS 101152-94-7

Name: Milnacipran hydrochloride CAS No.:101152-94-7 Molecular Structure: Molecular Structure of 101152-94-7 (Milnacipran hydrochloride) Formula: C15H23ClN2O Molecular Weight : 282.81 Synonyms: Cyclopropanecarboxamide,2-(aminomethyl)-N,N-diethyl-1-phenyl-, monohydrochloride,(57264057,1R,2S)-rel-(9CI);Milnacipran

N-Ethyl-N-benzylaniline-3'-sulfonic acid CAS 101-11-1

Name: Benzenesulfonic acid,3-[(ethylphenylamino)methyl]- CAS No.:101-11-1 Molecular Structure: Molecular Structure of 101-11-1 (Benzenesulfonic acid,3-[(ethylphenylamino)methyl]-) Formula: C15H17NO3S Molecular Weight : 291.36 Synonyms: m-Toluenesulfonicacid, a-(N-ethylanilino)-

Decyltrimethylammonium chloride CAS 10108-87-9

Name: 1-Decanaminium,N,N,N-trimethyl-, chloride (1:1) CAS No.:10108-87-9 Molecular Structure: Molecular Structure of 10108-87-9 (1-Decanaminium,N,N,N-trimethyl-, chloride (1:1)) Formula: C13H30N.Cl Molecular Weight : 235.84 Synonyms: 1-Decanaminium,N,N,N-trimethyl-, chloride (9CI);Ammonium, decyltrimethyl-, chloride

3-Carboxy-5-nitrophenylboronic acid CAS 101084-81-5

Name: Benzoicacid, 3-borono-5-nitro- CAS No.:101084-81-5 Molecular Structure: Molecular Structure of 101084-81-5 (Benzoicacid, 3-borono-5-nitro-) Formula: C7H6BNO6 Molecular Weight : 210.94 Synonyms: (3-

Carboxy-5-nitrophenyl)boronicacid;3-(Dihydroxyboryl)-5-nitrobenzoic acid;3-borono-5-nitrobenzoic acid; EINECS:

5-NITRO-1H-PYRROLO[2,3-B]PYRIDINE CAS 101083-92-5

Name: 1H-Pyrrolo[2,3-b]pyridine,5-nitro- CAS No.:101083-92-5 Molecular Structure: Molecular Structure of 101083-92-5 (1H-Pyrrolo[2,3-b]pyridine,5-nitro-) Formula: C7H5N3O2 Molecular Weight : 163.13 Synonyms: 5-Nitro-1H-pyrrolo[2,3-b]pyridine;1H-pyrrolo[2,3-b]pyridine, 5-nitro-;5-Nitro-1H-pyrrolo[2,3-b]pyridine;

2,6-DICHLOROPYRIDINE-3,4-DIAMINE CAS 101079-63-4

Name: 3,4-Pyridinediamine, 2,6-dichloro- CAS No.:101079-63-4 Molecular Structure: Molecular Structure of 101079-63-4 (3,4-Pyridinediamine, 2,6-dichloro-) Formula: C5H5Cl2N3 Molecular Weight : 176.99 Synonyms: 2,6-Dichloropyridine-3,4-diamine; EINECS: Density: 1.602 g/cm3 Melting Point: Boiling Point: 424.014 °C at 760

N,N-Dibenzylethanolamine CAS 101-06-4

Name: Ethanol,2-[bis(phenylmethyl)amino]- CAS No.:101-06-4 Molecular Structure: Molecular Structure of 101-06-4 (Ethanol,2-[bis(phenylmethyl)amino]-) Formula: C16H19NO Molecular Weight : 241.33 Synonyms: Ethanol,2-(dibenzylamino)- (6CI,7CI,8CI);2-(Dibenzylamino)ethanol;2-[Bis(phenylmethyl)amino]ethanol;D

phosphoric acid, copper salt CAS 10103-48-7

Name: Phosphoric acid, coppersalt (1:?) CAS No.:10103-48-7 Molecular Structure: Molecular Structure of 10103-48-7 (Phosphoric acid, coppersalt (1:?)) Formula: Cu. x H3 O4 P Molecular Weight : 380.580722 Synonyms: Phosphoricacid, copper salt (8CI,9CI); Copper orthophosphate; Copper phosphate; Primarycopper phosphate

Calcium phosphate CAS 10103-46-5

Name: Calcium phosphate CAS No.:10103-46-5 Molecular Structure: Molecular Structure of 10103-46-5 (Calcium phosphate) Formula: Ca. xH3O4P Molecular Weight : 502.31100 Synonyms: Phosphoricacid, calcium salt (8CI,9CI);Calcium phosphate;Crodax DP 30;Dikal 21;Dynafos;E 341;HAP 05NP;HAP 60NPG;KDV 15u;LF-CP-ZA;Man-Gill

N-(2-Ethyl-2-hexenylidene)-1-heneicosanamine CAS 101023-74-9

Name: 1-Heneicosanamine,N-(2-ethyl-2-hexen-1-ylidene)- CAS No.:101023-74-9 Molecular Structure: Molecular Structure of 101023-74-9 (1-Heneicosanamine,N-(2-ethyl-2-hexen-1-ylidene)-) Formula: C29H57 N Molecular Weight : 419.7696 Synonyms: 1-Heneicosanamine,N-(2-ethyl-2-hexenylidene)- (9CI); M 41; M 41 (corrosion

LITHIUM METASILICATE CAS 10102-24-6

Name: Silicic acid (H2SiO3),lithium salt (1:2) CAS No.:10102-24-6 Molecular Structure: Molecular Structure of 10102-24-6 (Silicic acid (H2SiO3),lithium salt (1:2)) Formula: Li2O3Si Molecular Weight : 89.97 Synonyms: Lithium silicate;Lithium silicon oxide(Li2SiO3);Lithiumsilicate (Li2SiO3) (6CI);Silicic acid (H2SiO3),

Sodium thiosulfate pentahydrate CAS 10102-17-7

Name: Sodium thiosulfate pentahydrate CAS No.:10102-17-7 Molecular Structure: Molecular Structure of 10102-17-7 (Sodium thiosulfate pentahydrate) Formula: H10Na2O8S2 Molecular Weight : 248.20 Synonyms: Sodiumthiosulfate pentahydrate;Sodium thiosulfate;Disodium thiosulfate pentahydrate;Thiosulfuric acid (H2S2O3),

SODIUM SULFITE HEPTAHYDRATE CAS 10102-15-5

Name: Sulfurous acid,disodium salt, heptahydrate (9CI) CAS No.:10102-15-5 Molecular Structure: Molecular Structure of 10102-15-5 (Sulfurous acid,disodium salt, heptahydrate (9CI)) Formula: H14Na2O10S Molecular Weight : 252.15 Synonyms: Sodiumsulfite heptahydrate; EINECS: 231-548-0 Density: Melting Point: Boiling

Palladium nitrate CAS 10102-05-3

Name: Palladium nitrate CAS No.:10102-05-3 Molecular Structure: Molecular Structure of 10102-05-3 (Palladium nitrate) Formula: Pd.(NO3)2 Molecular Weight : 230.4298 Synonyms: Nitricacid, palladium(2+) salt (8CI,9CI);Palladium nitrate (Pd(NO3)2) (6CI,7CI);Palladium dinitrate;Palladium(II) nitrate;Palladousnitrate;

Nickel sulfate hexahydrate CAS 10101-97-0

Name: Nickel sulfate hexahydrate CAS No.:10101-97-0 Molecular Structure: Molecular Structure of 10101-97-0 (Nickel sulfate hexahydrate) Formula: NiSO4.6(H2O) Molecular Weight : 262.89 Synonyms: Sulfuricacid, nickel(2+) salt (1:1), hexahydrate (8CI,9CI);Nickel monosulfatehexahydrate;Nickel sulfate (NiSO4)

1,2,3-TRIPHENYLGUANIDINE CAS 101-01-9

Name: Guanidine,N,N',N''-triphenyl- CAS No.:101-01-9 Molecular Structure: Molecular Structure of 101-01-9 (Guanidine,N,N',N''-triphenyl-) Formula: C19H17 N3 Molecular Weight : 287.39 Synonyms: Guanidine,1,2,3-triphenyl- (6CI,7CI,8CI); 1,2,3-Triphenylguanidine;N,N',N''-Triphenylguanidine; N,N',N''-Triphenylguanidine;

Lead(II) iodide CAS 10101-63-0

Name: Lead iodide (PbI2) CAS No.:10101-63-0 Molecular Structure: Molecular Structure of 10101-63-0 (Lead iodide (PbI2)) Formula: PbI2 Molecular Weight : 461.01 Synonyms: C.I. 77613;Lead diiodide;Lead diiodide (PbI2);Lead iodide;Plumbousiodide; EINECS: 233-256-9 Density: 6.16 g/mL at 25 °C(lit.) Melting Point: Boiling

Chromic sulfate CAS 10101-53-8

Name: Chromic sulfate CAS No.:10101-53-8 Molecular Structure: Molecular Structure of 10101-53-8 (Chromic sulfate) Formula: Cr2O12S3 Molecular Weight : 392.18 Synonyms: Baychrom A;Baychrom F;C.I. 77305;Chromic sulfate (Cr2(SO4)3);ChromitanB;Chromitan MS;Chromitan NA;Chromium III sulfate;Chromium sulfate;Chromiumsulfate

Lead(II) nitrate CAS 10099-74-8

Name: Lead dinitrate CAS No.:10099-74-8 Molecular Structure: Molecular Structure of 10099-74-8 (Lead dinitrate) Formula: N2O6Pb Molecular Weight : 331.21 Synonyms: Nitric acid,lead(2+) salt (8CI,9CI);Nitric acid, lead(2+)salt (2:1);Lead nitrate;Lead nitrate (Pb(NO3)2);Lead(2+) bis(nitrate);Lead(2+) nitrate;Lead(II)

99% CAS 10098-97-2

Name: Strontium, isotope ofmass 90 CAS No.:10098-97-2 Molecular Structure: Molecular Structure of 10098-97-2 (Strontium, isotope ofmass 90) Formula: Sr Molecular Weight : 89.9077 Synonyms: 90Sr; Sr90; Strontium-90 EINECS: Density: g/cm3 Melting Point: Boiling Point: °Cat760mmHg Flash Point: °C Solubility: Appearance:

L-Lysine hydrochloride CAS 10098-89-2

Name: L-Lysine hydrochloride CAS No.:10098-89-2 Molecular Structure: Molecular Structure of 10098-89-2 (L-Lysine hydrochloride) Formula: C6H14N2O2.xClH Molecular Weight : 182.64800 Synonyms: L-LYSINE HYDROCHLORIDE SOLUTION; (L)-lysine monohydrochloride; L-LYSINEHCL(FEED); L-Lysine/hydrochloric acid,(57264034,1:x);

4-nitrophenylglycolic acid CAS 10098-39-2

Name: Benzeneacetic acid, a-hydroxy-4-nitro- CAS No.:10098-39-2 Molecular Structure: Molecular Structure of 10098-39-2 (Benzeneacetic acid, a-hydroxy-4-nitro-) Formula: C8H7 N O5 Molecular Weight : 197.14488 Synonyms: Mandelicacid, p-nitro- (7CI,8CI); 4-Nitromandelic acid; 4-Nitrophenylglycolic acid; NSC126600;

5-[(3-chlorophenyl)amino]-Benzo[c]-2,6-naphthyridine-8-carboxylic acid CAS 1009820-21-6

Name: CX-4945 CAS No.:1009820-21-6 Molecular Structure: Molecular Structure of 1009820-21-6 (CX-4945) Formula: C19H12ClN3O2 Molecular Weight : 349.77 Synonyms: CX 4945;5-[(3-Chlorophenyl)amino]benzo[c]-2,6-naphthyridine-8-carboxylic acid; EINECS: Density: 1.491 g/cm3 Melting Point: Boiling Point: 568.45 °C at 760 mmHg

Benzyldimethylcarbinyl butyrate CAS 10094-34-5

Name: Butanoic acid,1,1-dimethyl-2-phenylethyl ester CAS No.:10094-34-5 Molecular Structure: Molecular Structure of 10094-34-5 (Butanoic acid,1,1-dimethyl-2-phenylethyl ester) Formula: C14H20O2 Molecular Weight : 220.34 Synonyms: Butyricacid, a,a-dimethylphenethyl ester (8CI);Dimethylbenzylcarbinyl

ATRACTYLOSIDE SODIUM SALT CAS 100938-11-2

Name: 19-Norkaur-16-en-18-oicacid, 15-hydroxy-2-[[2-O-(3-methyl-1-oxobutyl)-3,4-di-O-sulfo-b-D-glucopyranosyl]oxy]-, disodiumsalt,(57264028,2b,4a,15a)- (9CI) CAS No.:100938-11-2 Molecular Structure: Molecular Structure of 100938-11-2 (19-Norkaur-16-en-18-oicacid,

(2R,4R)-benzyl 4-hydroxy-2-(hydroxyMethyl)pyrrolidine-1-carboxylate CAS 1009335-39-0

0.99

3-[2,4-Bis((3S)-3-methylmorpholin-4-yl)pyrido[5,6-e]pyrimidin-7-yl]-N-methylbenzamide CAS 1009298-59-2

Name: AZD2014 CAS No.:1009298-59-2 Molecular Structure: Molecular Structure of 1009298-59-2 (AZD2014) Formula: C25H30N6O3 Molecular Weight : 462.54400 Synonyms: 3-[2,4-Bis((3S)-3-methylmorpholin-4-yl)pyrido[5,6-e]pyrimidin-7-yl]-N-methylbenzamide; EINECS: Density: 1.231 Melting Point: Boiling Point: Flash Point:

TYR-D-ALA-GLY-PHE-MET ACOH H2O CAS 100929-62-2

Name: L-Methionine,L-tyrosyl-D-alanylglycyl-L-phenylalanyl-, monoacetate (salt) (9CI) CAS No.:100929-62-2 Molecular Structure: Molecular Structure of 100929-62-2 (L-Methionine,L-tyrosyl-D-alanylglycyl-L-phenylalanyl-, monoacetate (salt) (9CI)) Formula: C28 H37 N5 O7 S Molecular Weight : 665.75 Synonyms:

1H-IMidazole, 5,5'-[1,1'-biphenyl]-4,4'-diylbis[2-(2S)-2-pyrrolidinyl-, hydrochloride (1:4) CAS 1009119-83-8

Name: 1H-IMidazole, 5,5'-[1,1'-biphenyl]-4,4'-diylbis[2-(2S)-2-pyrrolidinyl-, hydrochloride (1:4) CAS No.:1009119-83-8 Molecular Structure: Molecular Structure of 1009119-83-8 (1H-IMidazole, 5,5'-[1,1'-biphenyl]-4,4'-diylbis[2-(2S)-2-pyrrolidinyl-, hydrochloride (1:4)) Formula: Molecular Weight : 0 Synonyms:

99% CAS 1009-07-0

Name: Acetamide, N-(5-chloro-2-mercaptophenyl)- CAS No.:1009-07-0 Molecular Structure: Molecular Structure of 1009-07-0 (Acetamide, N-(5-chloro-2-mercaptophenyl)-) Formula: C8H8ClNOS Molecular Weight : Synonyms: Acetanilide, 5'-chloro-2'-mercapto- (7CI,8CI); EINECS: Density: Melting Point: Boiling Point: Flash Point:

1-(4-Pyridyl)piperazine CAS 1008-91-9

Name: 1-(4-Pyridyl)piperazine CAS No.:1008-91-9 Molecular Structure: Molecular Structure of 1008-91-9 (1-(4-Pyridyl)piperazine) Formula: C9H13N3 Molecular Weight : 163.2196 Synonyms: Piperazine,1-(4-pyridyl)-

GAMMA-PHENYL-GAMMA-BUTYROLACTONE CAS 1008-76-0

Name: 2(3H)-Furanone,dihydro-5-phenyl- CAS No.:1008-76-0 Molecular Structure: Molecular Structure of 1008-76-0 (2(3H)-Furanone,dihydro-5-phenyl-) Formula: C10H10 O2 Molecular Weight : 162.19 Synonyms: Butyricacid, 4-hydroxy-4-phenyl-, g-lactone (6CI,7CI); (RS)-g-Phenyl-g-butyrolactone;(?à)-g-Phenyl-g-butyrolactone;

4-methyl-N-{3-methyl-1-oxo-1-[(pyridin-2-ylmethyl)amino]butan-2-yl}cyclohexanecarboxamide CAS 10086-67-4

0.99

Benzyltrimethylammonium hydroxide CAS 100-85-6

Name: Benzyltrimethylammonium hydroxide CAS No.:100-85-6 Molecular Structure: Molecular Structure of 100-85-6 (Benzyltrimethylammonium hydroxide) Formula: C10H16N.HO Molecular Weight : 167.28 Synonyms: Benzenemethanaminium,N,N,N-trimethyl-, hydroxide (1:1);BTMAH 40;N,N,N-Trimethylbenzenemethanaminium

2-Methoxy-4-methylpyridine CAS 100848-70-2

Name: Pyridine,2-methoxy-4-methyl- CAS No.:100848-70-2 Molecular Structure: Molecular Structure of 100848-70-2 (Pyridine,2-methoxy-4-methyl-) Formula: C7H9NO Molecular Weight : 123.15 Synonyms:

FAPbBr3 CH(NH2)2PbBr3,Formamidinium Bromide Perovskite Formamidinium lead Bromide CAS 1008105-17-6

Name: FAPbBr3 CH(NH2)2PbBr3,Formamidinium Bromide Perovskite Formamidinium lead Bromide CAS No.:1008105-17-6 Molecular Structure: Molecular Structure of 1008105-17-6 (FAPbBr3 CH(NH2)2PbBr3,Formamidinium Bromide Perovskite Formamidinium lead Bromide) Formula: Molecular Weight : 0 Synonyms: FAPbBr3

Hydrocarbon oils, arom., mixed with polyethylene and polypropylene, pyrolyzed, light oil fraction CAS 100801-63-6

Name: Hydrocarbon oils, arom., mixed with polyethylene and polypropylene, pyrolyzed, light oil fraction CAS No.:100801-63-6 Molecular Structure: Molecular Structure of 100801-63-6 (Hydrocarbon oils, arom., mixed with polyethylene and polypropylene, pyrolyzed, light oil fraction) Formula: Unspecified Molecular Weight :

2-Anisidine-4-?-hydroxyethylsulfonesulfateester CAS 10079-20-6

Name: Ethanol,2-[(3-amino-4-methoxyphenyl)sulfonyl]-, 1-(hydrogen sulfate) CAS No.:10079-20-6 Molecular Structure: Molecular Structure of 10079-20-6 (Ethanol,2-[(3-amino-4-methoxyphenyl)sulfonyl]-, 1-(hydrogen sulfate)) Formula: C9H13NO7S2 Molecular Weight : 311.33 Synonyms: Ethanol,2-(4-methoxymetanilyl)-, hydrogen

1-Pyrrolidinecarboxylic acid, 2,2'-([1,1'-biphenyl]-4,4'-diyldi-1H-iMidazole-5,2-diyl)bis-, 1,1'-bis(1,1-diMethylethyl) ester,(57264009,2S,2'S)- CAS 1007882-23-6

Name: 1-Pyrrolidinecarboxylic acid, 2,2'-([1,1'-biphenyl]-4,4'-diyldi-1H-iMidazole-5,2-diyl)bis-, 1,1'-bis(1,1-diMethylethyl) ester,(57264010,2S,2'S)- CAS No.:1007882-23-6 Molecular Structure: Molecular Structure of 1007882-23-6 (1-Pyrrolidinecarboxylic acid,

(S)-2-[5-[4-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)phenyl]-1H-imidazol-2-y CAS 1007882-12-3

Name: (S)-2-[5-[4-(4,4,5,5-Tetramethyl-1,3,2-dioxaborolan-2-yl)phenyl]-1H-imidazol-2-y CAS No.:1007882-12-3 Molecular Structure: Molecular Structure of 1007882-12-3 ((S)-2-[5-[4-(4,4,5,5-Tetraethyl-1,3,2-dioxaborolan-2-yl)phenyl]-1H-imidazol-2-y) Formula: C24H34BN3O4 Molecular Weight : 439.36 Synonyms:

4,4'-methylenebis[3-hydroxy-2-naphthoic] acid, compound with 10,11-dihydro-N,N-dimethyl-5H-dibenz[b,f]azepine-5-propylamine (1:2) CAS 10075-24-8

Name: 2-Naphthalenecarboxylic acid, 4,4'-methylenebis[3-hydroxy-, compd. with 10,11-dihydro-N,N-dimethyl-5H-dibenz[b,f]azepine-5-propanamine (1:2) CAS No.:10075-24-8 Molecular Structure: Molecular Structure of 10075-24-8 (2-Naphthalenecarboxylic acid, 4,4'-methylenebis[3-hydroxy-, compd. with

2-Furanacetyl bromide, 5-methyl-alpha-oxo- (9CI) CAS 100750-53-6

Name: 2-Furanacetyl bromide,5-methyl-a-oxo- CAS No.:100750-53-6 Molecular Structure: Molecular Structure of 100750-53-6 (2-Furanacetyl bromide,5-methyl-a-oxo-) Formula: C7H5 Br O3 Molecular Weight : 217.0168 Synonyms: 2-Furanacetyl bromide, 5-methyl-alpha-oxo- (9CI) EINECS: Density: Melting Point: Boiling Point: Flash

4-(3,5-DIBROMO-2-PYRIDYLAZO)-N-ETHYL-N-(3-SULFOPROPYL)ANILINE, MONOSODIUM SALT, MONOHYDRATE CAS 100743-65-5

Name: 1-Propanesulfonic acid,3-[[4-[2-(3,5-dibromo-2-pyridinyl)diazenyl]phenyl]ethylamino]- CAS No.:100743-65-5 Molecular Structure: Molecular Structure of 100743-65-5 (1-Propanesulfonic acid,3-[[4-[2-(3,5-dibromo-2-pyridinyl)diazenyl]phenyl]ethylamino]-) Formula: C16H18 Br2 N4 O3 S Molecular Weight : 546.20923

[3-(MERCAPTOMETHYL)-2,4,6-TRIMETHYLPHENYL]METHANETHIOL CAS 10074-13-2

Name: 1,3-Benzenedimethanethiol,2,4,6-trimethyl- CAS No.:10074-13-2 Molecular Structure: Molecular Structure of 10074-13-2 (1,3-Benzenedimethanethiol,2,4,6-trimethyl-) Formula: C11H16 S2 Molecular Weight : 212.37 Synonyms: m-Xylene-a,a'-dithiol, 2,4,6-trimethyl- (7CI,8CI) EINECS: Density: 1.067g/cm3 Melting Point:

2-Formyl-3,4-dihydro-2H-pyran CAS 100-73-2

Name: 2H-Pyran-2-carboxaldehyde,3,4-dihydro- CAS No.:100-73-2 Molecular Structure: Molecular Structure of 100-73-2 (2H-Pyran-2-carboxaldehyde,3,4-dihydro-) Formula: C6H8O2 Molecular Weight : 112.14 Synonyms: 2-Formyl-3,4-dihydro-2H-pyran;2-Propenal,

N,N-DIMETHYLANTHRANILIC ACID METHYL ESTER CAS 10072-05-6

Name: Benzoic acid,2-(dimethylamino)-, methyl ester CAS No.:10072-05-6 Molecular Structure: Molecular Structure of 10072-05-6 (Benzoic acid,2-(dimethylamino)-, methyl ester) Formula: C10H13 N O2 Molecular Weight : 179.22 Synonyms: Anthranilic acid,N,N-dimethyl-, methyl ester (7CI,8CI); Methyl

5-Chloro-2-methylbenzothiazole CAS 1006-99-1

Name: Benzothiazole,5-chloro-2-methyl- CAS No.:1006-99-1 Molecular Structure: Molecular Structure of 1006-99-1 (Benzothiazole,5-chloro-2-methyl-) Formula: C8H6 Cl N S Molecular Weight : 183.66 Synonyms: 2-Methyl-5-chlorobenzothiazole;5-Chloro-2-methylbenzothiazole; NSC 8453 EINECS: 213-746-9 Density: 1.365g/cm3

2-Vinylpyridine CAS 100-69-6

Name: Pyridine,2-ethenyl- CAS No.:100-69-6 Molecular Structure: Molecular Structure of 100-69-6 (Pyridine,2-ethenyl-) Formula: C7H7N Molecular Weight : 105.15 Synonyms: Pyridine,2-vinyl- (8CI);2-Ethenylpyridine;2-Pyridylethylene;NSC18255;a-Vinylpyridine; EINECS: 202-879-8 Density: 0.968 g/cm3 Melting Point: Boiling

Thioanisole CAS 100-68-5

Name: Thioanisole CAS No.:100-68-5 Molecular Structure: Molecular Structure of 100-68-5 (Thioanisole) Formula: C7H8S Molecular Weight : 124.21 Synonyms: Sulfide,methyl phenyl (6CI,8CI);(Methylthio)benzene;1-Phenyl-1-thiaethane;Anisole,thio-;Methyl phenyl sulfide;Methylphenyl thioether;NSC 57916;Phenyl

METHYL 3-HYDROXY-5-ISOXAZOLECARBOXYLATE CAS 10068-07-2

Name: 5-Isoxazolecarboxylicacid, 2,3-dihydro-3-oxo-, methyl ester CAS No.:10068-07-2 Molecular Structure: Molecular Structure of 10068-07-2 (5-Isoxazolecarboxylicacid, 2,3-dihydro-3-oxo-, methyl ester) Formula: C5H5NO4 Molecular Weight : 143.10 Synonyms: 5-Isoxazolecarboxylicacid, 3-hydroxy-, methyl ester

Cyclohexanone oxime CAS 100-64-1

Name: Cyclohexanone oxime CAS No.:100-64-1 Molecular Structure: Molecular Structure of 100-64-1 (Cyclohexanone oxime) Formula: C6H11NO Molecular Weight : 113.16 Synonyms: (Hydroxyimino)cyclohexane;Antioxidant D;NSC 6300;OxiKhim-Styrol; EINECS: 202-874-0 Density: 1.1 g/cm3 Melting Point: Boiling Point: 208 °C at 760

(1S)-2-chloro-1-(3,4-difluorophenyl)-1-ethanol CAS 1006376-60-8

Name: (1S)-2-chloro-1-(3,4-difluorophenyl)-1-ethanol CAS No.:1006376-60-8 Molecular Structure: Molecular Structure of 1006376-60-8 ((1S)-2-chloro-1-(3,4-difluorophenyl)-1-ethanol) Formula: C8H7ClF2O Molecular Weight : 192.5903864 Synonyms:

9-Anthraceneboronic acid CAS 100622-34-2

Name: Boronic acid,B-9-anthracenyl- CAS No.:100622-34-2 Molecular Structure: Molecular Structure of 100622-34-2 (Boronic acid,B-9-anthracenyl-) Formula: C14H11BO2 Molecular Weight : 222.05 Synonyms: 9-Anthraceneboronicacid (6CI);Boronic acid, 9-anthracenyl- (9CI);9-Anthrylboronic acid; EINECS: -0 Density: 1.26 g/cm3

N,N'-Diethyl-1,3-propanediamine CAS 10061-68-4

Name: 1,3-Propanediamine,N1,N3-diethyl- CAS No.:10061-68-4 Molecular Structure: Molecular Structure of 10061-68-4 (1,3-Propanediamine,N1,N3-diethyl-) Formula: C7H18N2 Molecular Weight : 130.2312 Synonyms: 1,3-Propanediamine,N,N'-diethyl-

trans-1,3-Dichloropropene CAS 10061-02-6

Name: 1-Propene,1,3-dichloro-,(57263987,1E)- CAS No.:10061-02-6 Molecular Structure: Molecular Structure of 10061-02-6 (1-Propene,1,3-dichloro-,(57263988,1E)-) Formula: C3H4Cl2 Molecular Weight : 110.97 Synonyms: 1-Propene,1,3-dichloro-,(57263989,E)-;Propene, 1,3-dichloro-,(57263990,E)-

Chromic chloride hexahydrate CAS 10060-12-5

Name: Chromic chloride hexahydrate CAS No.:10060-12-5 Molecular Structure: Molecular Structure of 10060-12-5 (Chromic chloride hexahydrate) Formula: CrCl3.6(H2O) Molecular Weight : 266.45 Synonyms: Chromiumchloride (CrCl3), hexahydrate (8CI,9CI);Chromiumchloride hexahydrate;Chromium trichloride

PHENYLMAGNESIUM CHLORIDE CAS 100-59-4

Name: Magnesium,chlorophenyl- CAS No.:100-59-4 Molecular Structure: Molecular Structure of 100-59-4 (Magnesium,chlorophenyl-) Formula: C6H5ClMg Molecular Weight : 136.86 Synonyms: Magnesium,phenyl-, chloride (6CI);Phenylmagnesium chloride (6CI);Chlorophenylmagnesium; EINECS: 202-868-8 Density: 0.98 g/mL at 20 °C

porphycene CAS 100572-96-1

Name: 21,22,23,24-Tetraazapentacyclo[16.2.1.12,5.18,11.112,15]tetracosa-1(21),2,4,6,8(23),9,11,13,15,17,19-undecaene CAS No.:100572-96-1 Molecular Structure: Molecular Structure of 100572-96-1 (21,22,23,24-Tetraazapentacyclo[16.2.1.12,5.18,11.112,15]tetracosa-1(21),2,4,6,8(23),9,11,13,15,17,19-undecaene) Formula:

Tricyclo[3.3.2.02,8]deca-3,6,9-triene CAS 1005-51-2

Name: Tricyclo[3.3.2.02,8]deca-3,6,9-triene CAS No.:1005-51-2 Molecular Structure: Molecular Structure of 1005-51-2 (Tricyclo[3.3.2.02,8]deca-3,6,9-triene) Formula: C10H10 Molecular Weight : 0 Synonyms: Bullvalene;Bullvalene (C10H10); NSC 120527 EINECS: Density: 1.091g/cm3 Melting Point: Boiling Point: 211.3°C at 760

1-dodecyl-5-oxopyrrolidine-3-carboxylic acid CAS 10054-21-4

Name: 3-Pyrrolidinecarboxylicacid, 1-dodecyl-5-oxo- CAS No.:10054-21-4 Molecular Structure: Molecular Structure of 10054-21-4 (3-Pyrrolidinecarboxylicacid, 1-dodecyl-5-oxo-) Formula: C17H31NO3 Molecular Weight : 297.43294 Synonyms: 1-Dodecyl-5-oxopyrrolidine-3-carboxylicacid; EINECS: 233-182-7 Density: 1.021 g/cm3

hexaethynylbenzene CAS 100516-61-8

Name: Benzene, hexaethynyl- CAS No.:100516-61-8 Molecular Structure: Molecular Structure of 100516-61-8 (Benzene, hexaethynyl-) Formula: C18H6 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

3-(nitrooxy)propane-1-ol CAS 100502-66-7

Name: 1,3-Propanediol, mononitrate CAS No.:100502-66-7 Molecular Structure: Molecular Structure of 100502-66-7 (1,3-Propanediol, mononitrate) Formula: C3H7NO4 Molecular Weight : Synonyms: EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport

CHROMIUM(II) CHLORIDE CAS 10049-05-5

Name: Chromium chloride(CrCl2) CAS No.:10049-05-5 Molecular Structure: Molecular Structure of 10049-05-5 (Chromium chloride(CrCl2)) Formula: Cl2Cr Molecular Weight : 122.90 Synonyms: Chromiumdichloride;Chromium(II) chloride;Chromous chloride; EINECS: 233-163-3 Density: g/cm3 Melting Point: Boiling Point: °Cat760mmHg

CV-6209 CAS 100488-87-7

Name: Pyridinium,2-(2-acetyl-6-methoxy-3,9-dioxo-4,8-dioxa-2,10-diazaoctacos-1-yl)-1-ethyl-,chloride (1:1) CAS No.:100488-87-7 Molecular Structure: Molecular Structure of 100488-87-7 (Pyridinium,2-(2-acetyl-6-methoxy-3,9-dioxo-4,8-dioxa-2,10-diazaoctacos-1-yl)-1-ethyl-,chloride (1:1)) Formula: C34H60 N3 O6 . Cl

2-Mercapto-4-hydroxy-5,6-diaminopyrimidine CAS 1004-76-8

Name: 4(1H)-Pyrimidinone,5,6-diamino-2,3-dihydro-2-thioxo- CAS No.:1004-76-8 Molecular Structure: Molecular Structure of 1004-76-8 (4(1H)-Pyrimidinone,5,6-diamino-2,3-dihydro-2-thioxo-) Formula: C4H6N4OS Molecular Weight : 158.18 Synonyms: Uracil,5,6-diamino-2-thio-

THIOGLYCOLIC ACID N-BUTYL ESTER CAS 10047-28-6

Name: Butyl thioglycolate CAS No.:10047-28-6 Molecular Structure: Molecular Structure of 10047-28-6 (Butyl thioglycolate) Formula: C6H12O2S Molecular Weight : 148.22 Synonyms: Thioglycolic acid n-butyl ester;n-Butyl thioglycolate;Aceticacid, mercapto-, butyl ester (6CI,7CI,8CI,9CI);Butyl

Ammonium iron(II) sulfate CAS 10045-89-3

Name: Sulfuric acid, iron(2+)ammonium salt (2:1:2) CAS No.:10045-89-3 Molecular Structure: Molecular Structure of 10045-89-3 (Sulfuric acid, iron(2+)ammonium salt (2:1:2)) Formula: Fe.(NH4)2.(SO4)2 Molecular Weight : 284.05 Synonyms: Sulfuricacid, ammonium iron(2+) salt (2:2:1) (8CI,9CI);Ammonium ferrous

Ethyl 1-methylbutyl cyanoacetate CAS 100453-11-0

Name: Hexanoic acid,2-cyano-2-ethyl-3-methyl-, ethyl ester CAS No.:100453-11-0 Molecular Structure: Molecular Structure of 100453-11-0 (Hexanoic acid,2-cyano-2-ethyl-3-methyl-, ethyl ester) Formula: C12H21NO2 Molecular Weight : 211.30 Synonyms: 2-Cyano-2-ethyl-3-methylhexanoic acid ethyl ester; EINECS: Density: 0.944

2,N-Dimethyl-N-(3,3-diphenylpropyl)-1-amino-2-propanol CAS 100442-33-9

Name: 2,N-Dimethyl-N-(3,3-diphenylpropyl)-1-amino-2-propanol CAS No.:100442-33-9 Molecular Structure: Molecular Structure of 100442-33-9 (2,N-Dimethyl-N-(3,3-diphenylpropyl)-1-amino-2-propanol) Formula: C20H27NO Molecular Weight : 297.43 Synonyms: 1-[(3,3-Diphenylpropyl)methylamino]-2-methyl-2-propanol; EINECS:

ALUMINUM POTASSIUM SULFATE CAS 10043-67-1

Name: Potassium aluminum sulfate CAS No.:10043-67-1 Molecular Structure: Molecular Structure of 10043-67-1 (Potassium aluminum sulfate) Formula: Al.2H2O4S.K Molecular Weight : 258.21 Synonyms: Aluminumpotassium sulfate (AlK(SO4)2) (6CI);Alum;Alum potassium;Aluminium potassiumbis(sulfate);Aluminum potassium

1,3,5-TRIMETHYLBORAZINE CAS 1004-35-9

Name: Borazine,1,3,5-trimethyl- CAS No.:1004-35-9 Molecular Structure: Molecular Structure of 1004-35-9 (Borazine,1,3,5-trimethyl-) Formula: C3H12B3N3 Molecular Weight : 122.5805 Synonyms:

Boron nitride CAS 10043-11-5

Name: Boron nitride CAS No.:10043-11-5 Molecular Structure: Molecular Structure of 10043-11-5 (Boron nitride) Formula: BN Molecular Weight : 24.82 Synonyms: Elbor;BN 40SHP;Borazon;Denka boron nitride GP;Denka GP;Elbor LO 10B1-100;Elbor RM;Geksanit R; EINECS: 233-136-6 Density: 2.29 g/cm3 Melting Point: Boiling Point:

99% CAS 10043-01-0

0.99

Lercanidipine CAS 100427-26-7

Name: Lercanidipine CAS No.:100427-26-7 Molecular Structure: Molecular Structure of 100427-26-7 (Lercanidipine) Formula: C36H41N3O6 Molecular Weight : 611.73 Synonyms: (+-)-2-((3,3-Diphenylpropyl)methylamino)-1,1-dimethylethyl methyl

Sodium picosulfate CAS 10040-45-6

Name: Sodium picosulfate CAS No.:10040-45-6 Molecular Structure: Molecular Structure of 10040-45-6 (Sodium picosulfate) Formula: C18H13NNa2O8S2 Molecular Weight : 481.42 Synonyms: Phenol,4,4'-(2-pyridinylmethylene)bis-, bis(hydrogen sulfate) (ester), disodium salt(9CI);Phenol, 4,4'-(2-pyridylmethylene)di-,

Ceramides CAS 100403-19-8

Name: Ceramides CAS No.:100403-19-8 Molecular Structure: Molecular Structure of 100403-19-8 (Ceramides) Formula: Molecular Weight : 397.63488 Synonyms: CAPROYL SPHINGOSINE;CERAMIDE 1;CERAMIDE 2;CERAMIDE 3;CERAMIDE 4;CERAMIDE 5;CERAMIDE 1 A;CERAMIDE 6 II EINECS: 309-560-3 Density: Melting Point: Boiling Point: Flash

1003-73-2 CAS 1003-73-2

Name: Pyridine, 3-methyl-,1-oxide CAS No.:1003-73-2 Molecular Structure: Molecular Structure of 1003-73-2 (Pyridine, 3-methyl-,1-oxide) Formula: C6H7NO Molecular Weight : 109.14 Synonyms: 3-Picoline,1-oxide (6CI,8CI);3-Methylpyridine 1-oxide;3-Methylpyridine N-oxide;3-Picoline N-oxide;NSC 18254;b-Picoline

2-Hydroxy-5-methylpyridine CAS 1003-68-5

Name: 2(1H)-Pyridinone,5-methyl- CAS No.:1003-68-5 Molecular Structure: Molecular Structure of 1003-68-5 (2(1H)-Pyridinone,5-methyl-) Formula: C6H7NO Molecular Weight : 109.13 Synonyms:

N-ACETYL-ALPHA-D-GLUCOSAMINE CAS 10036-64-3

Name: 2-Acetamido-2-deoxy-alpha-D-glucose CAS No.:10036-64-3 Molecular Structure: Molecular Structure of 10036-64-3 (2-Acetamido-2-deoxy-alpha-D-glucose) Formula: C8H15NO6 Molecular Weight : 221.2078 Synonyms: Glucopyranose,2-acetamido-2-deoxy-, a-D-

PHENOL-OD CAS 1003-66-3

Name: Phenol-d CAS No.:1003-66-3 Molecular Structure: Molecular Structure of 1003-66-3 (Phenol-d) Formula: C6H5 D O Molecular Weight : 95.12 Synonyms: Phenol-O-d EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information: MSDS:View

Magnesium sulfate heptahydrate CAS 10034-99-8

Name: Magnesium sulfate CAS No.:10034-99-8 Molecular Structure: Molecular Structure of 10034-99-8 (Magnesium sulfate) Formula: MgSO4.7(H2O) Molecular Weight : 120.3676 Synonyms: Magnesium sulfate;Magnesium Sulfate Heptahydrate;Magnesium sulfat; EINECS: 231-298-2 Density: 2.66 g/cm3 Melting Point: Boiling Point: 330 °C

Calcium sulfate hemihydrate CAS 10034-76-1

Name: Sulfuric acid, calciumsalt, hydrate (2:2:1) CAS No.:10034-76-1 Molecular Structure: Molecular Structure of 10034-76-1 (Sulfuric acid, calciumsalt, hydrate (2:2:1)) Formula: Ca.H2O4S.1/2H2O Molecular Weight : 145.15 Synonyms: UNII-3RW091J48V;Sulfuric acid, calcium salt (1:1), hemihydrate;Sulfuric acid, calcium

bromoacetylcarnitine CAS 10034-25-0

Name: bromoacetylcarnitine CAS No.:10034-25-0 Molecular Structure: Molecular Structure of 10034-25-0 (bromoacetylcarnitine) Formula: Molecular Weight : 0 Synonyms: bromoacetylcarnitine EINECS: Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes: Transport Information:

1,3,4,6-Tetra-O-acetyl-a-D-glucosamineHCl CAS 10034-20-5

Name: b-D-Glucopyranose,2-amino-2-deoxy-, 1,3,4,6-tetraacetate, hydrochloride (1:1) CAS No.:10034-20-5 Molecular Structure: Molecular Structure of 10034-20-5 (b-D-Glucopyranose,2-amino-2-deoxy-, 1,3,4,6-tetraacetate, hydrochloride (1:1)) Formula: C14H21 N O9 . Cl H Molecular Weight : 0 Synonyms:

Calcium phosphate monobasic CAS 10031-30-8

Name: Calcium phosphate monobasic CAS No.:10031-30-8 Molecular Structure: Molecular Structure of 10031-30-8 (Calcium phosphate monobasic) Formula: H6CaO9P2 Molecular Weight : 252.07 Synonyms: Phosphoricacid, calcium salt (2:1), monohydrate (8CI,9CI);Acid calcium phosphatehydrate;Calcium bis(dihydrogen phosphate)

LEAD(II) BROMIDE CAS 10031-22-8

Name: Lead bromide (PbBr2) CAS No.:10031-22-8 Molecular Structure: Molecular Structure of 10031-22-8 (Lead bromide (PbBr2)) Formula: Br2Pb Molecular Weight : 367.00800 Synonyms: Leadbromide;Lead

dibromide;Plumbous bromide; EINECS: 233-084-4 Density: 6.66 Melting Point: Boiling Point: 916oC Flash Point: Solubility:

(R)-(+)-ALPHA-HYDROXYBENZENE-ACETONITRILE CAS 10020-96-9

Name: Benzeneacetonitrile, a-hydroxy-,(57263918,aR)- CAS No.:10020-96-9 Molecular Structure: Molecular Structure of 10020-96-9 (Benzeneacetonitrile, a-hydroxy-,(57263919,aR)-) Formula: C8H7NO Molecular Weight : 133.15 Synonyms: Benzeneacetonitrile,a-hydroxy-,(57263920,R)-;Mandelonitrile,(57263921,+)-

Hydrolyzed vegetable protein CAS 100209-45-8

Name: Protein hydrolyzates, vegetable CAS No.:100209-45-8 Molecular Structure: Molecular Structure of 100209-45-8 (Protein hydrolyzates, vegetable) Formula: Unspecified Molecular Weight : 0 Synonyms: Protein hydrolyzates, vegetable;HYDROLYZED VEGETABLE PROTEIN;Vegetable protein hydrolyzate;Casein acid hydrolysate

Terephthaloyl chloride CAS 100-20-9

Name: Terephthaloyl chloride CAS No.:100-20-9 Molecular Structure: Molecular Structure of 100-20-9 (Terephthaloyl chloride) Formula: C8H4Cl2O2 Molecular Weight : 203.02 Synonyms: Terephthaloylchloride (8Cl);1,4-Benzenedicarbonyl chloride;1,4-Bis(chlorocarbonyl)benzene;1,4-Di(chlorocarbonyl)benzene;Benzene

Aluminum, 2-(2-quinolinyl)-1H-indene-1,3(2H)-dione sulfo derivs. complexes CAS 100208-62-6

Name: Aluminum,2-(2-quinolinyl)-1H-indene-1,3(2H)-dione sulfo derivs. complexes CAS No.:100208-62-6 Molecular Structure: Molecular Structure of 100208-62-6 (Aluminum,2-(2-quinolinyl)-1H-indene-1,3(2H)-dione sulfo derivs. complexes) Formula: Unspecified Molecular Weight : 273.28500 Synonyms: Aluminum,

Molybdenum tetrakis(dimethylamide) CAS 100207-68-9

Name: Molybdenum,tetrakis(N-methylmethanaminato)-,(57263912,T-4)- CAS No.:100207-68-9 Molecular Structure: Molecular Structure of 100207-68-9 (Molybdenum,tetrakis(N-methylmethanaminato)-,(57263913,T-4)-) Formula: C8H24 Mo N4 Molecular Weight : 0 Synonyms: Methanamine,N-methyl-, molybdenum complex EINECS: Density:

S-[2-[3-[[4-[[[(2R,3S,4R,5R)-5-(6-aminopurin-9-yl)-4-hydroxy-3-phosphonooxyoxolan-2-yl]methoxy-hydroxyphosphoryl]oxy-hydroxyphosphoryl]oxy-2-hydroxy-3,3-dimethylbutanoyl]amino]propanoylamino]ethyl] (E)-oct-2-enethioate CAS 10018-94-7

Name: Coenzyme A,S-(2E)-2-octenoate CAS No.:10018-94-7 Molecular Structure: Molecular Structure of 10018-94-7 (Coenzyme A,S-(2E)-2-octenoate) Formula: C29H48 N7 O17 P3 S Molecular Weight : 0 Synonyms: Coenzyme A,S-2-octenoate,(57263909,E)- (8Cl); 2,3-trans-Octenoyl coenzyme A; Oct-2-trans-enoyl-CoA;Oct-trans-2-enoyl

CYCLOHEXYLSUCCINATE CAS 10018-78-7

Name: Butanedioic acid,1-cyclohexyl ester CAS No.:10018-78-7 Molecular Structure: Molecular Structure of 10018-78-7 (Butanedioic acid,1-cyclohexyl ester) Formula: C10H16 O4 Molecular Weight : 0 Synonyms: Butanedioicacid, monocyclohexyl ester (9Cl); Succinic acid, cyclohexyl ester (7Cl);Succinic acid, monocyclohexyl

1,4-DIISOPROPYLBENZENE CAS 100-18-5

Name: 1,4-Diisopropylbenzene CAS No.:100-18-5 Molecular Structure: Molecular Structure of 100-18-5 (1,4-Diisopropylbenzene) Formula: C12H18 Molecular Weight : 162.30 Synonyms: Benzene,p-diisopropyl- (8Cl); 1,4-Bis(1-methylethyl)benzene; 1,4-Bis(isopropyl)benzene;1,4-Diisopropylbenzene; NSC 84198;

4-(DIETHYLAMINO)BENZHYDRAZIDE CAS 100139-54-6

Name: Benzoic acid,4-(diethylamino)-, hydrazide CAS No.:100139-54-6 Molecular Structure: Molecular Structure of 100139-54-6 (Benzoic acid,4-(diethylamino)-, hydrazide) Formula: C11H17N3O Molecular Weight : 207.27 Synonyms: Benzoicacid, p-diethylamino-, hydrazide (6Cl);4-(Diethylamino)benzenecarbohydrazide; EINECS:

3,8-Dibromophenanthroline CAS 100125-12-0

Name: 1,10-Phenanthroline,3,8-dibromo- CAS No.:100125-12-0 Molecular Structure: Molecular Structure of 100125-12-0 (1,10-Phenanthroline,3,8-dibromo-) Formula: C12H6 Br2 N2 Molecular Weight : 340.01 Synonyms: 3,8-Dibromo-1,10-phenanthroline;3,8-Dibromophenanthroline EINECS: Density: 1.915 Melting Point: Boiling Point:

DIBENZOFURAN-4-BORONIC ACID CAS 100124-06-9

Name: Boronic acid,B-4-dibenzofuranyl- CAS No.:100124-06-9 Molecular Structure: Molecular Structure of 100124-06-9 (Boronic acid,B-4-dibenzofuranyl-) Formula: C12H9BO3 Molecular Weight : 212.01 Synonyms: 4-Dibenzofuranboronicacid (6CI);Boronic acid, 4-dibenzofuranyl-

3-METHYL-5-(TRIFLUOROMETHYL)PYRAZOLE CAS 10010-93-2

Name: 1H-Pyrazole,3-methyl-5-(trifluoromethyl)- CAS No.:10010-93-2 Molecular Structure: Molecular Structure of 10010-93-2 (1H-Pyrazole,3-methyl-5-(trifluoromethyl)-) Formula: C5H5F3N2 Molecular Weight : 150.10 Synonyms: Pyrazole,3(or 5)-methyl-5(or 3)-(trifluoromethyl)- (7CI);Pyrazole,3-methyl-5-(trifluoromethyl)-

4-Dimethylaminobenzaldehyde CAS 100-10-7

Name: 4-Dimethylaminobenzaldehyde CAS No.:100-10-7 Molecular Structure: Molecular Structure of 100-10-7 (4-Dimethylaminobenzaldehyde) Formula: C9H11NO Molecular Weight : 149.19 Synonyms: p-Dimethylaminobenzaldehyde;Benzaldehyde,p-(dimethylamino)-

3,4-DIHYDRO-1H-PYRIDO[1,2-A]PYRIMIDIN-6(2H)-ONE CAS 1000981-74-7

Name: 6H-Pyrido[1,2-a]pyrimidin-6-one, 1,2,3,4-tetrahydro- CAS No.:1000981-74-7 Molecular Structure: Molecular Structure of 1000981-74-7 (6H-Pyrido[1,2-a]pyrimidin-6-one, 1,2,3,4-tetrahydro-) Formula: C8 H10 N2 O Molecular Weight : 150.1778 Synonyms: 1,2,3,4-Tetrahydro-6H-pyrido[1,2-a]pyrimidin-6-one; EINECS: Density:

Glycerides, fish-oil CAS 100085-40-3

Name: Glycerides, fish-oil CAS No.:100085-40-3 Molecular Structure: Molecular Structure of 100085-40-3 (Glycerides, fish-oil) Formula: Unspecified Molecular Weight : 0 Synonyms: Fish-oilglycerides; EINECS: 309-181-3 Density: Melting Point: Boiling Point: Flash Point: Solubility: Appearance: Hazard Symbols: Risk Codes:

Aloe vera oil CAS 100084-89-7

Name: Aloe capensis, ext. CAS No.:100084-89-7 Molecular Structure: Molecular Structure of 100084-89-7 (Aloe capensis, ext.) Formula: Unspecified Molecular Weight : 0 Synonyms: ALOE VERA OIL;aloe oil;Aloe capensis, ext.;Einecs 309-126-3 EINECS: 309-126-3 Density: Melting Point: Boiling Point: Flash Point: Solubility:

N,N'-(2,2-dimethylpropylidene)hexamethylenediamine CAS 1000-78-8

Name: 1,6-Hexanediamine,N1,N6-bis(2,2-dimethylpropylidene)- CAS No.:1000-78-8 Molecular Structure: Molecular Structure of 1000-78-8 (1,6-Hexanediamine,N1,N6-bis(2,2-dimethylpropylidene)-) Formula: C16H32 N2 Molecular Weight : 0 Synonyms: 1,6-Hexanediamine,N,N'-bis(2,2-dimethylpropylidene)- (9CI);

4-Methoxybenzoyl chloride CAS 100-07-2

Name: Benzoylchloride, 4-methoxy- CAS No.:100-07-2 Molecular Structure: Molecular Structure of 100-07-2 (Benzoylchloride, 4-methoxy-) Formula: C8H7ClO2 Molecular Weight : 170.60 . Synonyms: p-Anisoylchloride (7CI,8CI);4-Anisoyl chloride;4-Methoxybenzoic acid chloride;4-Methoxyphenylchloroform;Anisoyl

Bis(trimethylsilyl)carbodiimide CAS 1000-70-0

Name: Silanamine,N,N'-methanetetraylbis[1,1,1-trimethyl- CAS No.:1000-70-0 Molecular Structure: Molecular Structure of 1000-70-0 (Silanamine,N,N'-methanetetraylbis[1,1,1-trimethyl-) Formula: C7H18 N2 Si2 Molecular Weight : 186.45 Synonyms: Carbodiimide,bis(trimethylsilyl)- (7CI,8CI);

7-tert-butyl 2-ethyl 3-bromo-5,6-dihydroimidazo[1,2-a]pyrazine-2,7(8H)-dicarboxylate CAS 1000576-75-9

Name: Imidazo[1,2-a]pyrazine-2,7(8H)-dicarboxylicacid, 3-bromo-5,6-dihydro-, 7-(1,1-dimethylethyl) 2-ethyl ester CAS No.:1000576-75-9 Molecular Structure: Molecular Structure of 1000576-75-9 (Imidazo[1,2-a]pyrazine-2,7(8H)-dicarboxylicacid, 3-bromo-5,6-dihydro-, 7-(1,1-dimethylethyl) 2-ethyl ester) Formula:

2-(6-methoxypyridin-2-yl)acetonitrile CAS 1000512-48-0

Name: 2-Pyridineacetonitrile,6-methoxy- CAS No.:1000512-48-0 Molecular Structure: Molecular Structure of 1000512-48-0 (2-Pyridineacetonitrile,6-methoxy-) Formula: C8H8N2O Molecular Weight : 148.16 Synonyms: 2-(6-methoxy-2-pyridyl)acetonitrile;(6-Methoxypyridin-2-yl)acetonitrile; EINECS: Density: 1.115 g/cm3 Melting

4-Isoxazolecarboxylicacid,5-methyl-,methylester(6CI,9CI) CAS 100047-54-9

Name: 4-Isoxazolecarboxylicacid, 5-methyl-, methyl ester CAS No.:100047-54-9 Molecular Structure: Molecular Structure of 100047-54-9 (4-Isoxazolecarboxylicacid, 5-methyl-, methyl ester) Formula: C6H7NO3 Molecular Weight : 141.12 Synonyms: Methyl5-methylisoxazole-4-carboxylate;methyl

4-Aminopyridine-2-carboxylic acid CAS 100047-36-7

Name: 2-Pyridinecarboxylicacid, 4-amino- CAS No.:100047-36-7 Molecular Structure: Molecular Structure of 100047-36-7 (2-Pyridinecarboxylicacid, 4-amino-) Formula: C6H6N2O2 Molecular Weight : 138.12 Synonyms: Picolinicacid, 4-amino- (6CI); EINECS: -0 Density: 1.417 g/cm3 Melting Point: Boiling Point: 446.3 °C at 760

4-BROMO-3-IODO-1H-PYRROLO[2,3-B]PYRIDINE CAS 1000340-34-0

Name: 1H-Pyrrolo[2,3-b]pyridine, 4-bromo-3-iodo- CAS No.:1000340-34-0 Molecular Structure: Molecular Structure of 1000340-34-0 (1H-Pyrrolo[2,3-b]pyridine, 4-bromo-3-iodo-) Formula: C7H4BrIN2 Molecular Weight : Synonyms: 4-bromo-3-iodo-1H-pyrrolo[2,3-b]pyridine; 1H-Pyrrolo[2,3-b]pyridine, 4-bromo-3-iodo- EINECS:

Foral 85 CAS 8050-31-5

Name: Foral 85 CAS No.:8050-31-5 Molecular Weight : 0 Synonyms: ENDERE S;ESTAR GUM 8 BG;GLYCERYL HYDROGENATED ROSINATE;STAYBELITE ESTER 10;STAYBELITE ESTER 5;foral85;Glycerolrosinresin136;Resinacidsandresinacids;Disproportionated rosin, glycerol ester;Ester gum;Glycerol ester of disproportionated rosin;Glycerol ester

Indinavir CAS 150378-17-9

Name: Indinavir CAS No.:150378-17-9 Formula: C36H47N5O4 Molecular Weight : 613.79 Synonyms: D-erythro-Pentonamide,2,3,5-trideoxy-N-(2,3-dihydro-2-hydroxy-1H-inden-1-yl)-5-[2-[[(1,1-dimethylethyl)amino]carbonyl]-4-(3-pyridinylmethyl)-1-piperazinyl]-2-(phenylmethyl)-,[1(1S,2R),5(S)]-;Compound

Iron silicide (FeSi) CAS 12022-95-6

Name: Iron silicide (FeSi) CAS No.:12022-95-6 Formula: FeSi Molecular Weight : 83.93 Synonyms: Ironmonosilicide; Iron silicide (Fe0.5Si0.5) Density: g/cm3 Melting Point: Boiling Point: °Cat760mmHg Flash Point: °C

1,2,4-Tributylphosphorotrithioate CAS 78-48-8

Name: 1,2,4-Tributylphosphorotrithioate CAS No.:78-48-8 Formula: C12H27 O P S3 Molecular Weight : 314.54 Synonyms: Butylphosphorotrithioate (6CI); B 1,776; Butiphos; Butyl phosphorotrithioate((BuS)3PO); Chemagro B 1776; DEF; DEF 6; DEF Defoliant; De-Green; Fos-Fall A;Fosfall; S,S,S-Tributyl phosphorotrithioate;

Benzene,1,2,4-trichloro-5-[(4-chlorophenyl)sulfonyl]- CAS 116-29-0

Name: Benzene,1,2,4-trichloro-5-[(4-chlorophenyl)sulfonyl]- CAS No.:116-29-0 Formula: C12H6Cl4O2S Molecular Weight : 356.05 Synonyms: Tetradiphon;V 18;p-Chlorophenyl2,4,5-trichlorophenyl sulfone;Sulfone,p-chlorophenyl 2,4,5-trichlorophenyl

Propanoic acid, 3-phosphono- CAS 5962-42-5

Name: Propanoic acid, 3-phosphono- CAS No.:5962-42-5 Formula: C3H7O5P Molecular Weight : 154.06 Synonyms: Propionicacid, 3-phosphono- (6CI,7CI,8CI);1-Carboxyethane-2-phosphonic acid;2-Carboxyethane-1-phosphonic acid;2-Carboxyethanephosphonic acid;2-Carboxyethylphosphonic acid;Carboxyethylphosphonic acid;Phosphonic

Avermectin A1a,25-cyclohexyl-4'-O-de(2,6-dideoxy-3-O-methyl-a-L-arabino-hexopyranosyl)-5-demethoxy-25-de(1-methylpropyl)-22,23-dihydro-5-(hydroxyimino)-(9CI) CAS 165108-07-6

Name: Avermectin A1a,25-cyclohexyl-4'-O-de(2,6-dideoxy-3-O-methyl-a-L-arabino-hexopyranosyl)-5-demethoxy-25-de(1-methylpropyl)-22,23-dihydro-5-(hydroxyimino)-(9CI) CAS No.:165108-07-6 Formula: C43H63NO11 Molecular Weight :

Aurate(1-),tetrachloro-, potassium (1:1),(57263849,SP-4-1)- CAS 13682-61-6

Name: Aurate(1-),tetrachloro-, potassium (1:1),(57263850,SP-4-1)- CAS No.:13682-61-6 Formula: AuCl4.K Molecular Weight : 377.87 Synonyms: Aurate(1-),tetrachloro-, potassium (8CI);Aurate(1-), tetrachloro-, potassium,(57263851,SP-4-1)-(9CI);Potassium chloroaurate(III) (6CI);Potassium tetrachloroaurate(III)(7CI);Gold

Orotic acid monohydrate CAS 50887-69-9

Name: Orotic acid monohydrate CAS No.:50887-69-9 Formula: C5H4N2O4.H2O Molecular Weight : 174.11 Synonyms: 2,6-Dioxo-1,2,3,6-tetrahydropyrimidine-4-carboxylic acid hydrate; EINECS: 200-619-8 Density: 1.642 g/cm3 Boiling Point: 431.3 °C at 760 mmHg Flash Point: 214.6 °C Appearance: white to off-white crystalline

5-Bromoimidazo[1,2-a]pyridine CAS 69214-09-1

Name: 5-Bromoimidazo[1,2-a]pyridine CAS No.:69214-09-1 Formula: C7H5BrN2 Molecular Weight : 197.03 Synonyms: 5-BROMO-IMIDAZO[1,2-A]PYRIDINE; Density: 1.697 g/cm3

4,4'-Bipyridinium,1,1'-dimethyl- CAS 4685-14-7

Name: 4,4'-Bipyridinium,1,1'-dimethyl- CAS No.:4685-14-7 Formula: C12H14N2 Molecular Weight : 186.28 Synonyms: 1,1'-Dimethyl-4,4'-bipyridinium;1,1'-Dimethyl-4,4'-bipyridinium cation;Dimethyl viologen;Methyl viologendication;Methyl viologen ion(2+);Methyl

ACYCLOVIR CAS 59277-89-3

Name: Acyclovir CAS No.:59277-89-3 Formula: C8H11N5O3 Molecular Weight : 225.20 Synonyms: Acyclovir EINECS: 261-685-1 Density: 1.77 g/cm3 Boiling Point: 595 °C at 760 mmHg Flash Point: 313.6 °C Solubility: Water: 0.7 mg/mL Appearance: White to light yellow crystal powder

Dibenz[b,e]oxepin-2-aceticacid, 11-[3-(dimethylamino)propylidene]-6,11-dihydro-,(57263841,11Z)- CAS 113806-05-6

Name: Dibenz[b,e]oxepin-2-aceticacid, 11-[3-(dimethylamino)propylidene]-6,11-dihydro-,(57263842,11Z)- CAS No.:113806-05-6 Formula: C21H23NO3 Molecular Weight : 337.4122 Synonyms: Dibenz[b,e]oxepin-2-aceticacid, 11-[3-(dimethylamino)propylidene]-6,11-dihydro-,(

3-Bromofuran CAS 22037-28-1

Name: 3-Bromofuran CAS No.:22037-28-1 Formula: C4H3BrO Molecular Weight : 146.97 . Synonyms:Furan, 3-bromo-; EINECS: 244-744-6 Density: 1.662 g/cm3 Boiling Point: 103 °C at 760 mmHg Flash Point: 3.3 °C Appearance: Colorless to light yellow liquid

Benzene, ethenyl-, polymer with 1,3-butadiene CAS 9003-55-8

Name: Benzene, ethenyl-, polymer with 1,3-butadiene CAS No.:9003-55-8 Formula: C12H14 Molecular Weight : 474.72 Synonyms:Polyco 2410;Polyco 2430;Patelacol HB 504;DL 893;Dow 460;SD 345 (polymer);EKCh 26A;Synthemul 96-023;L 8801;ABR 60; Density: 1.04 g/mL at 25oC Boiling Point: 145.2 °C at 760 mmHg Flash Point: 31.1

1,2-Ethanediamine,N1,N1,N2-trimethyl- CAS 142-25-6

Name: 1,2-Ethanediamine,N1,N1,N2-trimethyl- CAS No.:142-25-6 Formula: C5H14N2 Molecular Weight : 102.18 Synonyms: 1,2-Ethanediamine,N,N,N'-trimethyl- (9CI);Ethylenediamine, N,N,N'-trimethyl-

Benzenemethanaminium,N-dodecyl-ar-ethyl-N,N-dimethyl-, chloride (1:1) CAS 27479-28-3

Name: Benzenemethanaminium,N-dodecyl-ar-ethyl-N,N-dimethyl-, chloride (1:1) CAS No.:27479-28-3 Formula: C23H42N.Cl Molecular Weight : 368.04 Synonyms: Ammonium,dodecyl(ar-ethylbenzyl)dimethyl-, chloride (8CI);Benzenemethanaminium, N-dodecyl-ar-ethyl-N,N-dimethyl-,chloride (9CI);Dodecyldimethyl(ethylbenzyl)ammonium

Zinc sulphate CAS 7733-02-0

Name: Zinc sulphate CAS No.:7733-02-0 Formula: ZnSO4 Molecular Weight : 161.43 Synonyms: Zinc sulfate (1:1);Zinc sulfate (ZnSO4);Zinc vitriol;Zinc(II) sulfate;Zincaps;Zincate;Zinco;Zincomed;Zin Sulphate Mono/Hepta;BiolectraZink;Bonazen;Bufopto Zinc Sulfate;Complexonat;Honny Fresh

Tetracosane,2,6,10,15,19,23-hexamethyl- CAS 111-01-3

Name: Tetracosane,2,6,10,15,19,23-hexamethyl- CAS No.:111-01-3 Formula: C30H62 Molecular Weight : 422.92 Synonyms: Dodecahydrosqualene;Fitoderm;Mild Finish 20P;NSC 6851;Nikko Squalane;Perhydrosqualene;Phtyosqualan;Phytiane LS;Phytosqualan;Pripure 3759;Pripure 379;Pripure SQV

Chlordane CAS 57-74-9

Name: 4,7-Methano-1H-indene,1,2,4,5,6,7,8,8-octachloro-2,3,3a,4,7,7a-hexahydro- CAS No.:57-74-9 Formula: C10H6 Cl8 Molecular Weight : 409.7814 Synonyms: 4,7-Methanoindan,1,2,4,5,6,7,8,8-octachloro-3a,4,7,7a-tetrahydro-

2(3H)-Naphthalenone,4,4a,5,6,7,8-hexahydro-4,4a-dimethyl-6-(1-methylethenyl)-,(57263830,4R,4aS,6R)- CAS 4674-50-4

Name: 2(3H)-Naphthalenone,4,4a,5,6,7,8-hexahydro-4,4a-dimethyl-6-(1-methylethenyl)-,(57263831,4R,4aS,6R)- CAS No.:4674-50-4 Formula: C15H22O Molecular Weight :

5,9,13,17-Nonadecatetraen-2-one,6,10,14,18-tetramethyl- CAS 6809-52-5

Name: 5,9,13,17-Nonadecatetraen-2-one,6,10,14,18-tetramethyl- CAS No.:6809-52-5 Formula: C23H38O Molecular Weight : 330.55 Synonyms: Selbex;Teprenona;Teprenonum;UNII-S8S8451A4O;6,10,14,18-Tetramethyl-5,9,13,17-nonadecatetraen-2-one, mixture of (5E,9E,13E) and (5Z,9E,13E) isomers; Density: 0.872 g/cm3 Boiling Point:

1H-Benz[e]indolium,2-[7-[1,3-dihydro-1,1-dimethyl-3-(4-sulfobutyl)-2H-benz[e]indol-2-ylidene]-1,3,5-heptatrien-1-yl]-1,1-dimethyl-3-(4-sulfobutyl)-,inner salt, sodium salt (1:1) CAS 3599-32-4

Name: 1H-Benz[e]indolium,2-[7-[1,3-dihydro-1,1-dimethyl-3-(4-sulfobutyl)-2H-benz[e]indol-2-ylidene]-1,3,5-heptatrien-1-yl]-1,1-dimethyl-3-(4-sulfobutyl)-,inner salt, sodium salt (1:1) CAS No.:3599-32-4 Formula: C43H47N2NaO6S2 Molecular Weight :

Propanamide CAS 79-05-0

Name: Propanamide CAS No.:79-05-0 Formula: C3H7NO Molecular Weight : 73.11 Synonyms: Propionamide(6CI,8CI);NSC 38708;Propanimidic acid;Propionic acid amide;Propionic amide; EINECS: 201-172-1 Density: 0.926 g/cm3 Boiling Point: 222.2 °C at 760 mmHg Flash Point: 84.9 °C Solubility: SOLUBLE Appearance: white to slightly

α-1-ACETYLMETHADOL CAS 1477-40-3

Name: α-1-ACETYLMETHADOL CAS No.:1477-40-3 Formula: C23H31 N O2 Molecular Weight : 353.55 Synonyms: 3-Heptanol,6-(dimethylamino)-4,4-diphenyl-, acetate (ester),(57263792,3S,6S)-(-)-(8CI);Benzeneethanol, b-[(2S)-2-(dimethylamino)propyl]-a-ethyl-b-phenyl-, acetate (ester),(57263793,aS)- (9CI); Benzeneethanol,

1,10-Diaminodecane CAS 646-25-3

Name: 1,10-Diaminodecane CAS No.:646-25-3 Formula: C10H24 N2 Molecular Weight : 172.36 Synonyms: 1,10-Decamethylenediamine;1,10-Decylenediamine;1,10-Diamino-n-decane;1,10-Diaminodecane;1,6-Decanediamine;Decamethylenediamine;NCI 10400;NSC 10400; EINECS: 211-471-9 Density: 0.857 g/cm3 Boiling Point: 140 °C12 mm

Cholan-24-oic acid,3,12-dihydroxy-,(57263788,3a,5b,12a)- CAS 83-44-3

Name: Cholan-24-oic acid,3,12-dihydroxy-,(57263789,3a,5b,12a)- CAS No.:83-44-3 Formula: C24H40O4 Molecular Weight : 392.58 Synonyms: 5b-Cholan-24-oic acid, 3a,12a-dihydroxy- (8CI);17b-[1-Methyl-3-carboxypropyl]-etiocholane-3a,12a-diol;3a,12a-Dihydroxy-5b-cholan-24-oic acid;3a,12a-Dihydroxy-5b-cholanic

Quaternaryammonium compounds, dimethylditallow alkyl, chlorides CAS 68783-78-8

Name: Quaternaryammonium compounds, dimethylditallow alkyl, chlorides CAS No.:68783-78-8 Molecular Weight : 0 Synonyms: Adogen 442-110P;Adogen 470;Armosoft M;Arquad 2T;Arquad 2T75;Dimethylditallow alkylammoniumchlorides;Praepagen 3345;Prapagen3445;Prapagen S 3445;QuartaminCO 86;Quaternium 48; EINECS: 272-207-6

Lithium,[(trimethylsilyl)methyl]- CAS 1822-00-0

Name: Lithium,[(trimethylsilyl)methyl]- CAS No.:1822-00-0 Formula: C4H11LiSi Molecular Weight : 94.16 Synonyms: Silane,tetramethyl-, lithium complex;(Lithiomethyl)trimethylsilane;[(Trimethylsilyl)methyl]lithium; Density: 0.65 g/mL at 25 °C Melting Point: Boiling Point: 35-36 °C Flash Point: -49 °C

3-Methoxybenzenethiol CAS 15570-12-4

Name: 3-Methoxybenzenethiol CAS No.:15570-12-4 Formula: C7H8OS Molecular Weight : 140.20 Synonyms: Benzenethiol,m-methoxy- (6CI,7CI,8CI);(m-Methoxyphenyl)thiol;3-Mercaptoanisole;3-Methoxythiophenol;m-Methoxybenzenethiol;m-Methoxythiophenol; EINECS: 239-617-7 Density: 1.112 g/cm3 Boiling Point: 223.5 °C at 760

Tetradecanoic acid,sodium salt (1:1) CAS 822-12-8

Name: Tetradecanoic acid,sodium salt (1:1) CAS No.:822-12-8 Formula: C14H28 O2 . Na Molecular Weight : 251.41 Synonyms: Myristicacid, sodium salt (8CI); Tetradecanoic acid, sodium salt (9CI); Nonsoul MN 1;Sodium myristate; Sodium tetradecanoate EINECS: 212-487-9 Density: g/cm3 Boiling Point: 319.6 °C at 760 mmHg Flash

[1,3]Dioxolo[4,5-g]cinnoline-3-carboxylicacid, 1-ethyl-1,4-dihydro-4-oxo- CAS 28657-80-9

Name: [1,3]Dioxolo[4,5-g]cinnoline-3-carboxylicacid, 1-ethyl-1,4-dihydro-4-oxo- CAS No.:28657-80-9 Formula: C12H10 N2 O5 Molecular Weight : 262.24 Synonyms: 1-Ethyl-3-carboxy-6,7-methylenedioxy-4-cinnolone;Cinobac; Cinobactin; Cinoxacin; Compound 64716; NSC 304467; Noxigram; Uronorm Density: 1.64g/cm3 Boiling Point:

PPY 12 CAS 30604-81-0

Name: PPY 12 CAS No.:30604-81-0 Formula: C4H5N Molecular Weight : 67.0892 Synonyms: Density: 0.99g/cm3 Boiling Point: 129.8°Cat760mmHg Flash Point: 33.3°C

Devarda's Alloy CAS 2246186

Name: Devarda's Alloy CAS No.:8049-11-4 Formula: AlCuZn Molecular Weight : 155.92 Synonyms: Copper alloy,base,Cu,Al,Zn (Devarda's alloy); Appearance: grey solid

Formamide,N-[(4-amino-2-methyl-5-pyrimidinyl)methyl]-N-[4-hydroxy-1-methyl-2-[[(tetrahydro-2-furanyl)methyl]dithio]-1-buten-1-yl]- CAS 804-30-8

Name: Formamide,N-[(4-amino-2-methyl-5-pyrimidinyl)methyl]-N-[4-hydroxy-1-methyl-2-[[(tetrahydro-2-furanyl)methyl]dithio]-1-buten-1-yl]- CAS No.:804-30-8 Formula: C17H26N4O3S2 Molecular Weight :

2-Propenal, 2-methyl- CAS 78-85-3

Name: 2-Propenal, 2-methyl- CAS No.:78-85-3 Formula: C4H6O Molecular Weight : 70.09 Synonyms: Methacrylaldehyde(8CI);2-Formyl-1-propene;2-Methacrolein;2-Methyl-2-propen-1-al;2-Methyl-2-propenal;2-Methylacrolein;2-Methylenepropanal;2-Methylpropenal;Isobutenal;Methacrylic aldehyde;NSC

Benzaldehyde,2,4,6-trihydroxy- CAS 487-70-7

Name: Benzaldehyde,2,4,6-trihydroxy- CAS No.:487-70-7 Formula: C7H6O4 Molecular Weight : 154.12 Synonyms: Formylphloroglucinol;NSC 38610;Phloroglucinaldehyde; EINECS: 207-663-7 Density: 1.598 g/cm3 Boiling Point: 319.1 °C at 760 mmHg Flash Point: 161 °C Appearance: pale pink crystalline powder

Glycine,N-[2-[[2-(dodecylamino)ethyl]amino]ethyl]- CAS 6843-97-6

Name: Glycine,N-[2-[[2-(dodecylamino)ethyl]amino]ethyl]- CAS No.:6843-97-6 Formula: C18H39 N3 O2 Molecular Weight : 355.51868 Synonyms: Anon LG;Dodecylbis(aminoethyl)glycine; Dodicin; Lebon 15;N-Lauryl-N''-(carboxymethyl)diethylenetriamine; N-Lauryldiethylenetriaminoaceticacid;

Acetamide,N-[3-[bis[2-(acetyloxy)ethyl]amino]-4-methoxyphenyl]- CAS 23128-51-0

Name: Acetamide,N-[3-[bis[2-(acetyloxy)ethyl]amino]-4-methoxyphenyl]- CAS No.:23128-51-0 Formula: C17H24N2O6 Molecular Weight : 352.38 Synonyms: p-Acetanisidide,3'-[bis(2-hydroxyethyl)amino]-, diacetate (ester)

Chromium oxide (CrO2) CAS 12018-01-8

Name: Chromium oxide (CrO2) CAS No.:12018-01-8 Formula: CrO2 Molecular Weight : 83.9949 Synonyms: accroxc;accroxr;Chromium oxide (CrO2);chromium(iv)dioxide;chromiumdioxide;chromiumdioxide(cro2);chromiumoxide(cro2);CrO2 EINECS: 234-630-4 Density: 4.89 g/cm3 Solubility: Insoluble in water Appearance: black tetrahedral

2,3-Dihydrobenzofuran CAS 496-16-2

Name: 2,3-Dihydrobenzofuran CAS No.:496-16-2 Formula: C8H8O Molecular Weight : 120.15 Synonyms: Benzofuran,1,2-dihydro- (3CI);2,3-Dihydro-1-benzofuran;2,3-Dihydrobenzo[b]furan;Benzofuran,2,3-dihydro-;Benzodihydrofuran;Coumaran;Dihydrobenzofuran;Kumaran; EINECS: 207-817-3 Density: 1.096 g/cm3 Boiling Point: 188.2 °C at

2-Propanone,1-(2-fluorophenyl)- CAS 2836-82-0

Name: 2-Propanone,1-(2-fluorophenyl)- CAS No.:2836-82-0 Formula: C9H9FO Molecular Weight : 152.17 Synonyms: 2-Propanone,(57263772,o-fluorophenyl)- (7CI);2-Propanone, 1-(o-fluorophenyl)- (8CI);1-(o-Fluorophenyl)-2-propanone;3-(2-Fluorophenyl)-2-propanone; EINECS: 220-627-5 Density: 1.077 g/cm3 Boiling Point: 266.2 °C

Orazamide dihydrate CAS 2574-78-9

Name: Orazamide dihydrate CAS No.:2574-78-9 Formula: C9H10N6O5.2H2O Molecular Weight : 282.21 Synonyms: Aicorat dihydrate;1,2,3,6-Tetrahydro-2,6-dioxopyrimidine-4-carboxylic acid, compoundwith 5-amino-1H-imidazole-4-carboxamide (1:1) dihydrate;5-amino-3H-imidazole-4-carboxamide; 2,6-dioxo-3H-pyrimidine-4-carboxylic

dl-alpha-Tocopherol CAS 10191-41-0

Name: dl-alpha-Tocopherol CAS No.:10191-41-0 Formula: C29H50O2 Molecular Weight : 430.71 Synonyms: 2H-1-Benzopyran-6-ol,3,4-dihydro-2,5,7,8-tetramethyl-2-(4,8,12-trimethyltridecyl)-;6-Chromanol,2,5,7,8-tetramethyl-2-(4,8,12-trimethyltridecyl)-

Cupric Nitrate CAS 3251-23-8

Name: Cupric Nitrate CAS No.:3251-23-8 Formula: CuN2O6 Molecular Weight : 187.56 Synonyms: Nitricacid, copper(2+) salt (8CI,9CI);Claycop;Nitric acid, copper(2+)salt (2:1);Copper nitrate(Cu(NO3)2);Copper(2+) nitrate;Copper(II) nitrate;Cupric dinitrate;Cupricnitrate; EINECS: 221-838-5 Density: 1.00 g/mL at 20 °C Boiling

2(3H)-Benzofuranone,3-(methoxymethylene)- CAS 40800-90-6

Name: 2(3H)-Benzofuranone,3-(methoxymethylene)- CAS No.:40800-90-6 Formula: C10H8O3 Molecular Weight : 176.17 Synonyms: 3-(Methoxymethylene)-2(3H)-benzofuranone;3-(a-Methoxy)methylenebenzofuran-2(3H)-one;(3Z)-3-(Methoxymethylene)-1-benzofuran-2(3H)-one; Density: 1.358 g/cm3 Boiling Point: 331.1 °C at 760 mmHg Flash

Benzoyl chloride,2,3,4,5,6-pentafluoro- CAS 2251-50-5

Name: Benzoyl chloride,2,3,4,5,6-pentafluoro- CAS No.:2251-50-5 Formula: C7ClF5O Molecular Weight : 230.52 Synonyms: Benzoylchloride, pentafluoro- (6CI,7CI,8CI,9CI);2,3,4,5,6-Pentafluorobenzoic acidchloride;2,3,4,5,6-Pentafluorobenzoyl chloride;NSC 97002;Perfluorobenzoyl chloride; EINECS: 218-842-4 Density: 1.684

1-Hexadecene CAS 629-73-2

Name: 1-Hexadecene CAS No.:629-73-2 Formula: C16H32 Molecular Weight : 224.43 Synonyms: 1-Cetene;1-n-Hexadecene;Cetene;Dialen 16;LAO-C 16;Linealene 16;NSC 60602;NeoSolv6;Neodene 16;n-Hexadec-1-ene;a-Hexadecene; EINECS: 248-131-4 Density: 0.779 g/cm3 Boiling Point: 284.8 °C at 760 mmHg Flash Point: 132.2 °C Appearance:

Benzeneacetyl chloride,a-methoxy-a-(trifluoromethyl)-,(57263727,aS)- CAS 20445-33-4

Name: Benzeneacetyl chloride,a-methoxy-a-(trifluoromethyl)-,(57263728,aS)- CAS No.:20445-33-4 Formula: C10H8 Cl F3 O2 Molecular Weight : 252.62 Synonyms: Benzeneacetylchloride, a-methoxy-a-(trifluoromethyl)-,(57263729,S)-;Hydratropoyl chloride, b,b,b-trifluoro-a-methoxy-,(57263730,+)- (8CI);

3-Quinolinecarboxylicacid, 4-hydroxy-7-(phenylmethoxy)-, ethyl ester CAS 17825-15-9

Name: 3-Quinolinecarboxylicacid, 4-hydroxy-7-(phenylmethoxy)-, ethyl ester CAS No.:17825-15-9 Formula: C19H17 N O4 Molecular Weight : 323.34 Synonyms: 3-Quinolinecarboxylicacid, 7-(benzyloxy)-4-hydroxy-, ethyl ester (8CI) Density: 1.248g/cm3 Boiling Point: 491°C at 760 mmHg Flash Point: 250.8°C

Benzene, decyl- CAS 104-72-3

Name: Benzene, decyl- CAS No.:104-72-3 Formula: C16H26 Molecular Weight : 218.38 Synonyms: Decane,1-phenyl- (6CI,7CI,8CI); 1-Phenyl-n-decane; 1-Phenyldecane; Decylbenzene; NSC74191; Phenyldecane; n-Decylbenzene EINECS: 203-230-1 Density: 0.858g/cm3 Boiling Point: 293 °C(lit.) Flash Point: 126.8°C Appearance:

Pangamic acid, sodiumsalt CAS 77700-02-8

Name: Pangamic acid, sodiumsalt CAS No.:77700-02-8 Molecular Weight : 0 Synonyms: PANGAMIC ACID SODIUM;PANGAMIC ACID SODIUM SALT

Stibine, triphenyl- CAS 603-36-1

Name: Stibine, triphenyl- CAS No.:603-36-1 Formula: C18H15Sb Molecular Weight : 353.08 Synonyms: NSC 2844;Triphenylantimony;Triphenylstibane;Triphenylstibine;BRN 3608108;CCRIS 1300;Trifenylstibin; EINECS: 210-037-6 Density: 1.434 g/cm3 Boiling Point: 377 °C Flash Point: 109 °C Solubility: insoluble in

2,6-Diaminotoluene CAS 823-40-5

Name: 2,6-Diaminotoluene CAS No.:823-40-5 Formula: C7H10N2 Molecular Weight : 122.17 Synonyms: Toluene-2,6-diamine(7CI,8CI);1,3-Diamino-2-methylbenzene;2,6-Diamino-1-methylbenzene;2,6-Toluenediamine;2-Methyl-m-phenylenediamine; EINECS: 212-513-9 Density: 1.107 g/cm3 Boiling Point: 284.2 °C at 760 mmHg Flash Point:

Sphingosine 1-phosphate CAS 26993-30-6

Name: Sphingosine 1-phosphate CAS No.:26993-30-6 Formula: C18H38NO5P Molecular Weight : 379.47 Synonyms: 4-Octadecene-1,3-diol,2-amino-, 1-(dihydrogen phosphate),(57263718,E)-D-erythro- (8CI);4-Octadecene-1,3-diol,2-amino-, 1-(dihydrogen phosphate), [R-[R*,S*-(E)]]-;(4E)-2-Amino-3-hydroxy-4-octadecen-1-yl dihydrogen

1,4-Bis((2,3-epoxypropoxy)methyl)cyclohexane CAS 14228-73-0

Name: 1,4-Bis((2,3-epoxypropoxy)methyl)cyclohexane CAS No.:14228-73-0 Formula: C14H24 O4 Molecular Weight : 256.34 Synonyms: Cyclohexane,1,4-bis[(2,3-epoxypropoxy)methyl]- (8CI);1,4-Bis(glycidoxymethyl)cyclohexane;1,4-Bis(hydroxymethyl)cyclohexane diglycidyl

5-Nitrobenzimidazole nitrate CAS 27896-84-0

Name: 5-Nitrobenzimidazole nitrate CAS No.:27896-84-0 Formula: C7H5N3O2.HNO3 Molecular Weight : 226.15 Synonyms: 1H-Benzimidazole, 5-nitro-, mononitrate;Benzimidazole, 5-nitro-, mononitrate;nitric acid; 5-nitro-3H-benzoimidazole;6-nitro-1H-benzo[d]imidazole nitrate;5(6)-Nitrobenzimidazole nitrate;6-Nitrobenzimidazole

Ethanethiol,2-[(3-aminopropyl)amino]-, 1-(dihydrogen phosphate), hydrate (1:3) CAS 112901-68-5

Name: Ethanethiol,2-[(3-aminopropyl)amino]-, 1-(dihydrogen phosphate), hydrate (1:3) CAS No.:112901-68-5 Formula: C5H21N2O6PS Molecular Weight : 268.27 Synonyms: Anifostine trihydrate;Ethanethiol,2-[(3-aminopropyl)amino]-, dihydrogen phosphate (ester), trihydrate (9CI); Density: 1.367g/cm3 Boiling Point: 441.7 °C at

3,5-Pyridinedicarboxylicacid, 1,4-dihydro-2,6-dimethyl-4-(3-nitrophenyl)-, 3-methyl5-[2-[methyl(phenylmethyl)amino]ethyl] ester, hydrochloride (1:1) CAS 54527-84-3

Name: 3,5-Pyridinedicarboxylicacid, 1,4-dihydro-2,6-dimethyl-4-(3-nitrophenyl)-, 3-methyl5-[2-[methyl(phenylmethyl)amino]ethyl] ester, hydrochloride (1:1) CAS No.:54527-84-3 Formula: C26H30ClN3O6 Molecular Weight : 515.9859 Synonyms: 3,5-Pyridinedicarboxylicacid, 1,4-dihydro-2,6-dimethyl-4-(3-nitrophenyl)-,

Iodic acid (HIO3),potassium salt (2:1) CAS 13455-24-8

Name: Iodic acid (HIO3),potassium salt (2:1) CAS No.:13455-24-8 Formula: HIO3.1/2K Molecular Weight : 390.91 Synonyms: Potassiumiodate (7CI);Hydrogen potassium iodate;Potassium hydrogen iodate (1:1:2);Potassium biiodat; EINECS: 236-650-9

Platinum,trimethyl[(1,2,3,4,5-h)-1-methyl-2,4-cyclopentadien-1-yl]- CAS 94442-22-5

Name: Platinum,trimethyl[(1,2,3,4,5-h)-1-methyl-2,4-cyclopentadien-1-yl]- CAS No.:94442-22-5 Formula: C9H16 Pt Molecular Weight : 319.30 Synonyms: Platinum,trimethyl(methylcyclopentadienyl)- (7CI); 1,3-Cyclopentadiene, 1-methyl-,platinum complex; (Methylcyclopentadienyl)trimethylplatinum;

Dodecanamine CAS 124-22-1

Name: Dodecanamine CAS No.:124-22-1 Formula: C12H27N Molecular Weight : 185.40 Synonyms: Lauramine;Laurinamine;Laurylamine;Monododecylamine;Monolaurylamine;Nissan Amine BB;n-Dodecylamine;n-Laurylamine;1-Aminododecane;1-Dodecylamine;Alamine 4;Amine BB;Armeen 12D;Farmin20D;HAN 12;Kemamine P 690; EINECS:

Imidazo[2,1-b]quinazolin-2(3H)-one,6,7-dichloro-1,5-dihydro-, hydrochloride (1:1) CAS 58579-51-4

Name: Imidazo[2,1-b]quinazolin-2(3H)-one,6,7-dichloro-1,5-dihydro-, hydrochloride (1:1) CAS No.:58579-51-4 Formula: C10H8Cl3N3O Molecular Weight : 292.55 Synonyms: Imidazo[2,1-b]quinazolin-2(3H)-one,6,7-dichloro-1,5-dihydro-, monohydrochloride (9CI);6,7-Dichloro-1,5-dihydroimidazo[2,1-b]quinazolin-2(3H)-one

Glycoproteins,bovine-whey CAS 84082-51-9

Name: Glycoproteins,bovine-whey CAS No.:84082-51-9 Molecular Weight : 0 Synonyms: Glycoproteins, bovine-whey

5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid,7-[[(2Z)-2-(2-amino-4-thiazolyl)-2-(methoxyimino)acetyl]amino]-3-[(1Z)-2-(4-methyl-5-thiazolyl)ethenyl]-8-oxo-,(57263703,2,2-dimethyl-1-oxopropoxy)methyl ester,(57263704,6R,7R)- CAS 117467-28-4

Name: 5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid,7-[[(2Z)-2-(2-amino-4-thiazolyl)-2-(methoxyimino)acetyl]amino]-3-[(1Z)-2-(4-methyl-5-thiazolyl)ethenyl]-8-oxo-,(57263705,2,2-dimethyl-1-oxopropoxy)methyl ester,(57263706,6R,7R)- CAS No.:117467-28-4 Formula: C25H28N6O7S3 Molecular Weight :

Calcium sulfide (CaS) CAS 20548-54-3

Name: Calcium sulfide (CaS) CAS No.:20548-54-3 Formula: CaS Molecular Weight : 72.14 Synonyms: Calciumsulfide (8CI); Baumol; C.I. 77245; C.I. Pigment White 9; Calcium monosulfide;Pigment White 9 Density: g/cm3 Boiling Point: °Cat760mmHg Flash Point: °C

1H-Imidazole-5-propanamide,N-[(1S,2R,3S)-1-(cyclohexylmethyl)-3-cyclopropyl-2,3-dihydroxypropyl]-a-[[(2S)-2-[[(1,1-dimethylethyl)sulfonyl]methyl]-1-oxo-3-phenylpropyl]amino]-,(57263697,aS)- CAS 126222-34-2

Name: 1H-Imidazole-5-propanamide,N-[(1S,2R,3S)-1-(cyclohexylmethyl)-3-cyclopropyl-2,3-dihydroxypropyl]-a-[[(2S)-2-[[(1,1-dimethylethyl)sulfonyl]methyl]-1-oxo-3-phenylpropyl]amino]-,(57263698,aS)- CAS No.:126222-34-2 Formula: C33H50 N4 O6 S Molecular Weight :

1-Dodecylpyridinium bromide CAS 104-73-4

Name: 1-Dodecylpyridinium bromide CAS No.:104-73-4 Formula: C17H30BrN Molecular Weight : 328.33 Synonyms: 1-Dodecylpyridiniumbromide (6CI);Pyridinium, 1-dodecyl-, bromide (8CI,9CI);Dodecylpyridinium bromide;Isothan Q 4;Laurosept;Laurylpyridinium bromide;N-Dodecylpyridinium bromide;N-Laurylpyridinium bromide; EINECS:

4-Cyclohexene-1,2,3-triol,6-amino-4-(hydroxymethyl)-,(57263693,1S,2S,3R,6S)- CAS 38231-86-6

Name: 4-Cyclohexene-1,2,3-triol,6-amino-4-(hydroxymethyl)-,(57263694,1S,2S,3R,6S)- CAS No.:38231-86-6 Formula: C7H13NO4 Molecular Weight : 175.18 Synonyms: (+)-Valienamine; Density: 1.525 g/cm3 Boiling Point: 377.3 °C at 760 mmHg Flash Point: 182 °C Appearance: Slightly yellow solid

1,3-Benzenedicarboxamide,5-(acetylamino)-N1,N3-bis(2,3-dihydroxypropyl)-2,4,6-triiodo- CAS 31127-80-7

Name: 1,3-Benzenedicarboxamide,5-(acetylamino)-N1,N3-bis(2,3-dihydroxypropyl)-2,4,6-triiodo- CAS No.:31127-80-7 Formula: C16H20I3N3O7 Molecular Weight : 747.06 Synonyms: 1,3-Benzenedicarboxamide,5-(acetylamino)-N,N'-bis(2,3-dihydroxypropyl)-2,4,6-triiodo- (9CI);Isophthalamide,

3-Quinolinecarboxylicacid,7-(3-amino-1-pyrrolidinyl)-8-chloro-1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxo-,hydrochloride (1:1) CAS 105956-99-8

Name: 3-Quinolinecarboxylicacid,7-(3-amino-1-pyrrolidinyl)-8-chloro-1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxo-,hydrochloride (1:1) CAS No.:105956-99-8 Formula: C17H17ClFN3O3.HCl Molecular Weight :

b-D-Galactopyranoside,(57263688,3b,4a,16a)-16,23,28-trihydroxyoleana-11,13(18)-dien-3-yl 6-deoxy-3-O-b-D-glucopyranosyl- CAS 58316-41-9

Name: b-D-Galactopyranoside,(57263689,3b,4a,16a)-16,23,28-trihydroxyoleana-11,13(18)-dien-3-yl 6-deoxy-3-O-b-D-glucopyranosyl- CAS No.:58316-41-9 Formula: C42H68O13 Molecular Weight : 780.99 Synonyms: SaikosaponinB2; Density: 1.355 g/cm3 Boiling Point: 909.877 °C at 760 mmHg Flash Point: 504.075 °C

Poly(oxy-1,2-ethanediyl),a-isooctadecyl-w-hydroxy- CAS 52292-17-8

Name: Poly(oxy-1,2-ethanediyl),a-isooctadecyl-w-hydroxy- CAS No.:52292-17-8 Formula: (C2H4 O)n C18 H38 O Molecular Weight : 0 Synonyms: Aerosurf66 E10; Agnique MMF; Aldol 66E20; Arosurf 66E10; Arosurf 66E2; Arosurf 66E20;Arosurf 90E2; Arosurf AA 66E20; Arosurf ISA 20E; Arosurf MSF; Emalex 1810;Emalex 1815; Emalex

4,5,6,7-Tetrachloro-2',4',5',7'-tetraiodofluorescein dipotassium salt CAS 11121-48-5

Name: 4,5,6,7-Tetrachloro-2',4',5',7'-tetraiodofluorescein dipotassium salt CAS No.:11121-48-5 Formula: C20H2Cl4I4Na2O5 Molecular Weight : 1017.64 Synonyms: C.I.AcidRed94(45440);redno.105;rose;RB;TETRACHLOROTETRAIODOFLUORESCEIN SODIUM SALT;TETRAIODOTETRACHLOROFLUORESCEIN DISODIUM SALT;SODIUM

Phosphoric acid,7-chlorobicyclo[3.2.0]hepta-2,6-dien-6-yl dimethyl ester CAS 23560-59-0

Name: Phosphoric acid,7-chlorobicyclo[3.2.0]hepta-2,6-dien-6-yl dimethyl ester CAS No.:23560-59-0 Formula: C9H12 Cl O4 P Molecular Weight : 250.63 Synonyms: Bicyclo[3.2.0]heptane,phosphoric acid deriv.; 4-(O,O-

Dimethylphosphoryl)-5-chlorobicyclo[3.2.0]hepta-1,4-diene;Heptenofos; Heptenophos; Hostaquick; Ragadan; XOE

Percoll CAS 65455-52-9

Name: Percoll CAS No.:65455-52-9 Molecular Weight : 0 Synonyms: PERCOLL;PERCOLL(R);PERCOLL(TM);CENTRICOLL;CENTRICOLL(TM);PERCOLL, PLANT CELL CULTURE TESTED;PERCOLL CELL CULTURE TESTED;silica particles coated with polyvinylpyrrolidone Density: 1.1 g/mL at 25 °C(lit.)

Magnesate(2-),bis[L-aspartato(2-)-kN,kO1]-, hydrogen (1:2),(57263678,T-4)- CAS 18962-61-3

Name: Magnesate(2-),bis[L-aspartato(2-)-kN,kO1]-, hydrogen (1:2),(57263679,T-4)- CAS No.:18962-61-3 Formula: C8H12MgN2O8 Molecular Weight : 288.49 Synonyms: Magnesate(2-), bis[L-aspartato(2-)-N,O1]-,dihydrogen,(57263680,T-4)-;Magnesate(2-), bis[L-aspartato(2-)-kN,kO1]-, dihydrogen,(57263681,T-4)-(9CI);L-Aspartic

1-Hexadecen-3-ol,3,7,11,15-tetramethyl- CAS 505-32-8

Name: 1-Hexadecen-3-ol,3,7,11,15-tetramethyl- CAS No.:505-32-8 Formula: C20H40O Molecular Weight : 296.54 Synonyms:NSC 93744; EINECS: 208-008-8 Density: 0.851 g/cm3 Boiling Point: 327.8 °C at 760 mmHg Flash Point: 143.5 °C Solubility: Practically insoluble in water Appearance: clear to yellow oily liquidi

Cadmium nitrate tetrahydrate CAS 10022-68-1

Name: Cadmium nitrate tetrahydrate CAS No.:10022-68-1 Formula: N2O6.Cd.4H2O Molecular Weight : 308.50 Synonyms: Nitricacid, cadmium salt, tetrahydrate (8CI,9CI);C.I. 77192;Cadmium dinitratetetrahydrate;Cadmium nitrate (Cd(NO3)2) tetrahydrate;Nitric acid, cadmiumsalt, hydrate (2:1:4); EINECS: 232-146-8 Density: 2.45

KANAMYCIN ACID SULFATE CAS 64013-70-3

Name: KANAMYCIN ACID SULFATE CAS No.:64013-70-3 Molecular Weight : 680.66 Synonyms: KANAMYCIN ACID SULPHATE;KANAMYCIN A ACID SULFATE;KANAMYCIN ACID SULFATE;Kanamycindisulfate;kanamycin acid sulfate from*streptomyces kanamyce;Kanamycin Sulfate(Sterile);KANAMYCIN ACID SULFATE FROM*STREPTOMYCES KANAMYCETI;KANAMYCIN ACID

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Acetamide,N-(phenylmethyl)- CAS 588-46-5

Name: Acetamide,N-(phenylmethyl)- CAS No.:588-46-5 Formula: C9H11 N O Molecular Weight : 149.19 Synonyms: Acetamide,N-benzyl- (6CI,7CI,8CI); N-(Phenylmethyl)acetamide; N-Acetylbenzylamine;N-Benzylacetamide; NSC 8065 Density: 1.031g/cm3 Boiling Point: 157 °C / 2mmHg Flash Point: 194.5°C

Cyclohexane,1,4-bis(isocyanatomethyl)- CAS 10347-54-3

Name: Cyclohexane,1,4-bis(isocyanatomethyl)- CAS No.:10347-54-3 Formula: C10H14 N2 O2 Molecular Weight : 194.23036 Synonyms: Isocyanicacid, 1,4-cyclohexylenedimethylene ester (7CI);1,4-Bis(isocyanatomethyl)cyclohexane Density: 1.14g/cm3 Boiling Point: 294.4°Cat760mmHg Flash Point: 119.5°C

Tallow, hydrogenated CAS 2239276

Name: Tallow, hydrogenated CAS No.:8030-12-4 Molecular Weight : 0 Synonyms: Beeftallow, hydrogenated;Carolac;Hydrofol Glyceride T 57N;Hydrogenated beeftallow;Hydrogenated tallow;T 57N; EINECS: 232-442-7

ZIRCONIUM-TITANIUM ALLOY CAS 50646-37-2

Name: ZIRCONIUM-TITANIUM ALLOY CAS No.:50646-37-2 Molecular Weight : 139.09 Synonyms: ZIRCONIUM-TITANIUM ALLOY;Zirconium-titanium alloy, -100 mesh, 50:50 atomic %;Alloy-100mesh,50:50atomic%;Zirconium Titanium powder, -100 mesh

4,6(1H,5H)-Pyrimidinedione,5-ethyldihydro-5-(1-methylbutyl)-2-thioxo-, sodium salt (1:1) CAS 71-73-8

Name: 4,6(1H,5H)-Pyrimidinedione,5-ethyldihydro-5-(1-methylbutyl)-2-thioxo-, sodium salt (1:1) CAS No.:71-73-8 Formula: C11H18 N2 O2 S . Na Molecular Weight : 265.36 Synonyms: 4,6(1H,5H)-Pyrimidinedione,5-ethyldihydro-5-(1-methylbutyl)-2-thioxo-, monosodium salt (9CI);Barbituricacid, 5-ethyl-5-(1-methylbutyl)-2-thio-,

Ouabain CAS 630-60-4

Name: Ouabain CAS No.:630-60-4 Formula: C29H44O12 Molecular Weight : 584.65 Synonyms: Ouabain(8CI);Acocantherin;Astrobain;Gratibain;Gratus strophanthin;Kombetin;NSC

D(+)-Glucose CAS 50-99-7

Name: D(+)-Glucose CAS No.:50-99-7 Formula: C6H12O6 Molecular Weight : 180.16 Synonyms: (+)-Glucose;Glucose;C*Dry GL 01934;CPC hydrate;Cartose;Cerelose;Cerelose2001;Clearsweet 95;Clintose L;Corn sugar;Dextropur;Dextrose;Dextrosol;Glucodin;Glucolin;D-Glucose;Glucosteril;Goldsugar;Grape sugar;Hi-Mesh;Maxim Energy

Androst-5-ene-3,17-diol,(57263662,3b,17b)- CAS 521-17-5

Name: Androst-5-ene-3,17-diol,(57263663,3b,17b)- CAS No.:521-17-5 Formula: C19H30O2 Molecular Weight : 290.44 Synonyms: Androst-5-ene-3b,17b-diol (7CI,8CI);3b,17b-Androst-5-enediol;3b,17b-Dihydroxyandrost-5-ene;Androst-5-enediol;Hermaphrodiol;NSC 12163;D5-Androstene-3b,17b-diol;D5-Androstenediol; EINECS:

Acid Blue 9 CAS 2650-18-2

Name: Acid Blue 9 CAS No.:2650-18-2 Formula: C37H34N2Na2O9S3 Molecular Weight : 783.01 Synonyms: A.F. Blue No. 1;AF Blue No. 1;Acid Blue;AE-SF;Acilan Turquoise Blue AE;Alphazurine FGND;Bleu Brilliant FCF;C.I. 42090;C.I. Acid Blue 9, diammonium salt;C.I. Direct Brown 78, Diammonium Salt;CI 42090;CI Acid Blue 9,

Acid Blue 277 CAS 25797-81-3

Name: Acid Blue 277 CAS No.:25797-81-3 Formula: C24H22N3NaO8S2 Molecular Weight : 567.57 Synonyms: Best Acid Brilliant Blue 4R;C.I. 61203;C.I. Acid Blue 277;Erio Blue BGL;Nylosan Blue E 4R;Taitilon Blue GLF-T;Tectilon Blue 4R;Triacid Fast Blue 4RL; EINECS: 247-269-2

Benzoic acid,4-[2-(1-piperidinyl)ethoxy]-, hydrochloride (1:1) CAS 84449-80-9

Name: Benzoic acid,4-[2-(1-piperidinyl)ethoxy]-, hydrochloride (1:1) CAS No.:84449-80-9 Formula: C14H20ClNO3 Molecular Weight : 285.77 Synonyms: Benzoicacid, 4-[2-(1-piperidinyl)ethoxy]-, hydrochloride (9CI);4-(2-Piperidinoethoxy)benzoic acid hydrochloride; EINECS: 282-882-9 Boiling Point: 413.2 °C at 760 mmHg Flash

Benzenemethanamine,2,4-difluoro- CAS 72235-52-0

Name: Benzenemethanamine,2,4-difluoro- CAS No.:72235-52-0 Formula: C7H7F2N Molecular Weight : 143.13 Synonyms: (2,4-Difluorophenyl)methanamine;2,4-Difluorobenzenemethanamine;[(2,4-Difluorophenyl)methyl]amine; EINECS: 276-502-0 Density: 1.214 g/cm3 Boiling Point: 177.8 °C at 760 mmHg Flash Point: 68.3 °C Appearance:

Methyl 3-phenylpropionate CAS 103-25-3

Name: Methyl 3-phenylpropionate CAS No.:103-25-3 Formula: C10H12O2 Molecular Weight : 164.20 Synonyms: Methyldihydrocinnamate;Methyl hydrocinnamate;Methyl b-phenylpropionate;NSC 10128;NSC 404188;b-Phenylpropionic acid methylester;Hydrocinnamicacid, methyl ester (6CI,7CI,8CI);3-Phenylpropanoic acid methyl

Allylestrenol CAS 432-60-0

Name: Allylestrenol CAS No.:432-60-0 Formula: C21H32O Molecular Weight : 300.53 Synonyms: Estr-4-en-17-ol,17-(2-propenyl)-,(57263653,17b)- (9CI);Estr-4-en-17b-ol, 17-allyl- (6CI,8CI);17a-Allyl-17b-hydroxy-D4-estren;17a-Allyl-4-oestren-17b-ol;17a-Allylestr-4-en-17b-ol;17a-Allylestrenol;Gestanin;Gestanol;Gestanon;NSC

- ## Benzenemethanol,a-(1-aminoethyl)-2,5-dimethoxy-,hydrochloride (1:1) CAS 61-16-5

  Name: Benzenemethanol,a-(1-aminoethyl)-2,5-dimethoxy-,hydrochloride (1:1) CAS No.:61-16-5 Formula: C11H17NO3.ClH Molecular Weight : 247.72 Synonyms: Benzenemethanol,a-(1-aminoethyl)-2,5-dimethoxy-,hydrochloride (9CI);Methoxamine hydrochloride

- ## Tamoxifen citrate CAS 54965-24-1

  Name: Tamoxifen citrate CAS No.:54965-24-1 Formula: C26H29NO.C6H8O7 Molecular Weight : 563.70 Synonyms: Farmifeno;Oncomox;Tamoxifen citrate (JAN/USP);Tafoxen;Kessar;Tamoxasta;Tamoxifenum [INN-Latin];Nolvadex (TN);Soltamox;Noxitem;Oncotam;Ethanamine, 2-(4-(1,2-diphenyl-1-butenyl)phenoxy)-N,N-dimethyl,(57263649,Z)-,

- ## Neoandrographolide CAS 27215-14-1

  Name: Neoandrographolide CAS No.:27215-14-1 Formula: C26H40O8 Molecular Weight : 480.59 Synonyms: {(1R,4aS,8aS)-1,4a-dimethyl-6-methylidene-5-[2-(2-oxo-2,5-dihydrofuran-3-yl)ethyl]decahydronaphthalen-1-yl}methyl beta-D-glucopyranoside;2(5H)-Furanone,

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Fudosteine CAS 13189-98-5

Name: Fudosteine CAS No.:13189-98-5 Formula: C6H13NO3S Molecular Weight : 179.26 Synonyms: SS 320A;S-(3-Hydroxypropyl)cysteine;(2R)-2-amino-3-(3-hydroxypropylsulfanyl)propanoic acid;Cleanal;Cleanal (TN);L-Cysteine, S-(3-hydroxypropyl)-;(-)-3-((3-Hydroxypropyl)thio)-L-alanine;Alanine, 3-((3-hydroxypropyl)thio)-, L-

- ## L-Alpha-Glycerophosphocholine CAS 28319-77-9

Name: Ethanaminium,2-[[[(2R)-2,3-dihydroxypropoxy]hydroxyphosphinyl]oxy]-N,N,N-trimethyl-, innersalt CAS No.:28319-77-9 Formula: C8H20NO6P Molecular Weight : 257.22 Synonyms: Alfoscerate de choline;BRN 6062450;Calcium (glycerophosphate de);Choline alfoscerate;Choline alphoscerate;Cholini alfosceras;Cholini

- ## L-Theanine CAS 3081-61-6

Name: L-Glutamine,N-ethyl- CAS No.:3081-61-6 Formula: C7H14N2O3 Molecular Weight : 174.2 Synonyms: Glutamine,N-ethyl-, L- (6CI,7CI,8CI);NSC 21308;Ng-Ethyl-L-glutamine;Suntheanine;Theanin;Theanine;L-Theanine (Ngamma-ethyl-L-glutamine); EINECS: 221-379-0 Density: 1.171 g/cm3 Boiling Point: 430.2 °C at 760 mmHg Flash

- ## Methylsulfonylmethane CAS 67-71-0

Name: Methyl sulfone CAS No.:67-71-0 Formula: C2H6O2S Molecular Weight : 94.13 Synonyms: Methane,sulfonylbis- (9CI);Methyl sulfone (6CI,8CI);Dimethyl sulfone;Dimethylsulphone;Lignisul MSM;MSM;Methylsulfonylmethane;NSC 63345;Opti MSM;Opti-dimethyl sulfone;Dimethyl sulfone (MSM); EINECS: 200-665-9 Density: 1.139

N-Acetyl-L-Tryptophan CAS 1218-34-4

Name: N-Acetyl-L-tryptophan CAS No.:1218-34-4 Formula: C13H14N2O3 Molecular Weight : 246.29 Synonyms: Ac-Trp-OH;Ac-L-Trp-OH;Tryptophan,N-acetyl-, L- (8CI);(S)-N-Acetyltryptophan;Acetyl-L-tryptophan;Acetyltryptophan;L-N-Acetyltryptophan;N-Acetyltryptophan;NSC 90726; EINECS: 214-935-9 Density: 1.33g/cm3 Boiling Point:

N-Acetyl-DL-Phenylalanine CAS 2901-75-9

Name: Phenylalanine,N-acetyl- CAS No.:2901-75-9 Formula: C11H13NO3 Molecular Weight : 207.23 Synonyms: Alanine,N-acetyl-3-phenyl-, DL- (8CI);DL-Phenylalanine, N-acetyl-;Afalanine;DL-2-Acetamido-3-phenylpropionic acid;DL-N-Acetyl-3-phenylalanine;DL-N-Acetylphenylalanine;N-Acetyl-3-phenyl-DL-alanine;NSC 43242; EINECS:

Benzaldehyde,3-methoxy-4-methyl- CAS 24973-22-6

Name: Benzaldehyde,3-methoxy-4-methyl- CAS No.:24973-22-6 Formula: C9H10O2 Molecular Weight : 150.17 Synonyms: m-Anisaldehyde,4-methyl- (7CI,8CI);3-Methoxy-p-tolualdehyde; Density: 1.062 g/cm3 Boiling Point: 246.2 °C at 760 mmHg Flash Point: 105.7 °C

1,4-Bis(2-cyanostyryl)benzene CAS 13001-39-3

Name: 1,4-Bis(2-cyanostyryl)benzene CAS No.:13001-39-3 Formula: C24H16N2 Molecular Weight : 332.40 Synonyms: Benzonitrile,2,2'-(p-phenylenedivinylene)di- (7CI,8CI);2-[(E)-2-[4-[(E)-2-(2-cyanophenyl)ethenyl]phenyl]ethenyl]benzonitrile; EINECS: 235-835-1 Density: 1.18 g/cm3 Boiling Point: 575.7 °C at 760 mmHg Flash

- ## 3-Pyrrolidinol,hydrochloride (1:1),(57263631,3R)- CAS 104706-47-0

Name: 3-Pyrrolidinol,hydrochloride (1:1),(57263632,3R)- CAS No.:104706-47-0 Formula: C4H9NO.HCl Molecular Weight : 123.58 Synonyms: R-3-Hydroxypyrrolidine Hcl; (R)-(-)-3-Pyrrolidinol Hydrochloride; (R)-3-Pyrrolidinol hydrochloride; (3R)-pyrrolidin-3-ol,hydrochloride; (R)-3-Hydroxypyrrolidine hydrochloride;

- ## 2-Pentanol CAS 6032-29-7

Name: 2-Pentanol CAS No.:6032-29-7 Formula: C5H12O Molecular Weight : 88.15 Synonyms: 1-Methyl-1-butanol;1-Methylbutanol;2-Hydroxypentane;2-Pentyl alcohol;DL-2-Pentanol;Methylpropylcarbinol;sec-Amyl alcohol;sec-Pentyl alcohol; EINECS: 227-907-6 Density: 0.809 g/cm3 Boiling Point: 118.8 °C at 760 mmHg Flash Point: 33.9

- ## 1-Naphthalenesulfonicacid, 4-[2-(2-hydroxy-1-naphthalenyl)diazenyl]-, sodium salt (1:1) CAS 1658-56-6

Name: 1-Naphthalenesulfonicacid, 4-[2-(2-hydroxy-1-naphthalenyl)diazenyl]-, sodium salt (1:1) CAS No.:1658-56-6 Formula: C20H14N2O4S.Na Molecular Weight : 401.38 Synonyms: Naphthalene Red J (6CI);2-Naphthol

RedJ;Acid Leather Red ROC;Acid Red 88;AcidRed A;Acid Red AV;Acid Red G;Acid Rose AV;Amacid Fast Red A;Anadurm

Oxiconazole nitrate CAS 64211-46-7

Name: Oxiconazole nitrate CAS No.:64211-46-7 Formula: C18H13Cl4N3O.HNO3 Molecular Weight : 492.14 Synonyms: (1Z)-N-[(2,4-dichlorobenzyl)oxy]-1-(2,4-dichlorophenyl)-2-(1H-imidazol-1-yl)ethanimine nitrate;(1Z)-N-[(2,4-dichlorobenzyl)oxy]-1-(2,4-dichlorophényl)-2-(1H-imidazol-1-yl)éthanimine

2,3-Dichlorobenzoylcyanide CAS 77668-42-9

Name: 2,3-Dichlorobenzoylcyanide CAS No.:77668-42-9 Formula: C8H3Cl2NO Molecular Weight : 200.02 Synonyms: (2,3-Dichlorophenyl)oxoacetonitrile;2,3-Dichlorobenzoyl nitrile; EINECS: 278-747-9 Density: 1.446 g/cm3 Boiling Point: 309.4 °C at 760 mmHg Flash Point: 140.9 °C

Rhein CAS 478-43-3

Name: Rhein CAS No.:478-43-3 Formula: C15H8O6 Molecular Weight : 284.22 . Synonyms: Monorhein;4,5-DiOH-anthraquinone-2-COOH;Chrysazin-3-carboxylic acid;Rhein (1,8-dihydroxy-3-carboxyl anthraquinone);4,5-Dihydroxyanthraquinone-2-carboxylic acid;Rheic acid;9,10-Dihydro-4,5-dihydroxy-9,10-dioxo-2-anthracene carboxylic

frage CAS 112199-06-1

Name: L-Arginine,

Benzamide,4-amino-5-bromo-N-[2-(diethylamino)ethyl]-2-methoxy- CAS 4093-35-0

Name: Benzamide,4-amino-5-bromo-N-[2-(diethylamino)ethyl]-2-methoxy- CAS No.:4093-35-0 Formula: C14H22 Br N3 O2 Molecular Weight : 344.30 Synonyms: o-Anisamide,4-amino-5-bromo-N-[2-(diethylamino)ethyl]- (7CI,8CI); 2-Methoxy-4-amino-5-bromo-N,N-diethylaminoethylbenzamide;Bromopride; Emepride;

Diacerein CAS 13739-02-1

Name: Diacerein CAS No.:13739-02-1 Formula: C19H12O8 Molecular Weight : 368.29 Synonyms: 2-Anthracenecarboxylic acid, 4,5-bis(acetyloxy)-9,10-dihydro-9,10-dioxo- (9CI);4,5-diacetyloxy-9,10-dioxo-anthracene-2-carboxylic acid;4,5-Bis(acetyloxy)-9,10-dihydro-9,10-dioxo-2-anthracenecarboxylic

Bezafibrate CAS 41859-67-0

Name: Bezafibrate CAS No.:41859-67-0 Formula: C19H20ClNO4 Molecular Weight : 361.82 Synonyms: Bezatol SR;LO 44;2-(p-(2-(p-Chlorobenzamido)ethyl)phenoxy)-2-methylpropionic acid;Bezalip;Prestwick_724;BF-759;Cedur;2-[4-[2-[(4-chlorobenzoyl)amino]ethyl]phenoxy]-2-methyl-propanoic

Glufosinate-ammonium CAS 77182-82-2

Name: Glufosinate-ammonium CAS No.:77182-82-2 Formula: C5H15N2O4P Molecular Weight : 198.19 Synonyms: Butanoicacid, 2-amino-4-(hydroxymethylphosphinyl)-, monoammonium salt (9CI);Ammoniumglufosinate;Basta;Basta 150SL;Basta Fl;Basta LS;Buster;Dash;Finale;Finale 14SL;Glufosinate monoammonium salt; EINECS:

1H-Perimidin-2-amine,hydrobromide, hydrate (1:1:1) CAS 313223-13-1

Name: 1H-Perimidin-2-amine,hydrobromide, hydrate (1:1:1) CAS No.:313223-13-1 Formula: C11H9 N3 . Br H . H2 O Molecular Weight : 264.12 Synonyms: 1H-Perimidin-2-amine,monohydrobromide, monohydrate (9CI) Boiling Point: 421.7°C at 760 mmHg Flash Point: 208.8°C

2-(2-Methoxyphenoxy)ethylamine hydrochloride CAS 64464-07-9

Name: 2-(2-Methoxyphenoxy)ethylamine hydrochloride CAS No.:64464-07-9 Formula: C9H13NO2.HCl Molecular Weight : 203.67 Synonyms: Ethanamine,2-(2-methoxyphenoxy)-, hydrochloride (9CI);Ethylamine, 2-(o-methoxyphenoxy)-,hydrochloride (7CI); Boiling Point: 261.1 °C at 760 mmHg Flash Point: 120.7 °C Appearance: white

3-Fluorotoluene CAS 352-70-5

Name: 3-Fluorotoluene CAS No.:352-70-5 Formula: C7H7F Molecular Weight : 110.13 Synonyms: Toluene, m-fluoro- (8CI);m-Fluorotoluene;Toluene, m-fluoro-;1-chloro-4-fluoro-benzene;1-Fluoro-3-methylbenzene;1-Methyl-3-fluorobenzene;Benzene, 1-fluoro-3-methyl-;m-Fluorotoluene [UN2388] [Flammable liquid]; EINECS:

Sodium diethyldithiocarbamate CAS 148-18-5

Name: Sodium diethyldithiocarbamate CAS No.:148-18-5 Formula: C5H10NNaS2 Molecular Weight : 171.27 Synonyms: Sodium N,N-diethyldithiocarbamate;Carbamic acid, diethyldithio-, sodium salt;DEDC;Soxinol ESL;Diethyldithiocarbamic acid, sodium;Usaf ek-2596;DEDK;Ditiocarb sodium;Diethyldithiocarbamic acid, sodium

Dimethylaminopropanol CAS 3179-63-3

Name: Dimethylaminopropanol CAS No.:3179-63-3 Formula: C5H13NO Molecular Weight : 103.16282 [g/mol] Synonyms: 1-Propanol,

Dibenz[b,f][1,4]oxazepine,2-chloro-11-(1-piperazinyl)- CAS 14028-44-5

Name: Dibenz[b,f][1,4]oxazepine,2-chloro-11-(1-piperazinyl)- CAS No.:14028-44-5 Formula: C17H16ClN3O Molecular Weight : 313.79 Synonyms: 2-Chloro-11-(1-piperazinyl)dibenz[b,f][1,4]oxazepine; EINECS: 237-867-1 Density: 1.37 g/cm3 Boiling Point: 469.9 °C at 760 mmHg Flash Point: 238 °C Appearance: Crystalline solid

1-Acetylindoline CAS 16078-30-1

Name: 1-Acetylindoline CAS No.:16078-30-1 Formula: C10H11NO Molecular Weight : 161.20 Synonyms: 1H-Indole,1-acetyl-2,3-dihydro- (9CI);Indoline, 1-acetyl- (6CI,7CI,8CI);1-(2,3-Dihydroindol-1-yl)ethanone;1-(Indolin-1-yl)ethanone;1-Acetyl-2,3-dihydro-1H-indole;N-Acetylindoline; Density: 1.144 g/cm3 Boiling Point: 355.1

Bepotastine besilate CAS 190786-44-8

Name: Bepotastine besilate CAS No.:190786-44-8 Formula: C21H25ClN2O3.C6H6O3S Molecular Weight : 547.07 Synonyms: (+)-(S)-4-(4-((4-Chlorophenyl)(2-pyridyl)methoxy)piperidino)butyric acid monobenzenesulfonate;benzenesulfonic acid; 4-[4-[(S)-(4-chlorophenyl)-pyridin-2-yl-methoxy]-1-piperidyl]butanoic acid;Bepotastine

Acetic acid, iron(2+)salt (2:1) CAS 3094-87-9

Name: Acetic acid, iron(2+)salt (2:1) CAS No.:3094-87-9 Formula: C2H4O2.1/2Fe Molecular Weight : 87.97 Synonyms: Aceticacid, iron(2+) salt (8CI,9CI);Iron acetate [Fe(OAc)2];Irondiacetate;Iron(2+) acetate;Iron(II) acetate;Iron di(acetate); EINECS: 221-441-7 Boiling Point: 117.1 °C at 760 mmHg Flash Point: 40

Silanediol,1,1-diphenyl- CAS 947-42-2

Name: Silanediol,1,1-diphenyl- CAS No.:947-42-2 Formula: C12H12O2Si Molecular Weight : 216.31 Synonyms: Silanediol,diphenyl- (8CI,9CI);D 6150;Dihydroxydiphenylsilane;LS 4320;NSC 12561;SX 11;SX 11 (diol);TSL 8162; EINECS: 213-427-4 Density: 1.19 g/cm3 Boiling Point: 352.6 °C at 760 mmHg Flash Point: 167.1

5,5'-Diallyl-2,2'-biphenyldiol CAS 528-43-8

Name: 5,5'-Diallyl-2,2'-biphenyldiol CAS No.:528-43-8 Formula: C18H18O2 Molecular Weight : 266.36 Synonyms: 2,2'-Bichavicol(6CI);2,2'-Biphenyldiol, 5,5'-diallyl- (8CI);[1,1'-Biphenyl]-2,2'-diol,5,5'-di-2-propenyl- (9CI); Density: 1.107 g/cm3 Boiling Point: 401 °C at 760 mmHg Flash Point: 184.5 °C

SULPHADOXINE CAS 2447-57-6

Name: Sulfadoxine CAS No.:2447-57-6 Formula: C12H14N4O4S Molecular Weight : 310.33 Synonyms: Benzenesulfonamide,4-amino-N-(5,6-dimethoxy-4-pyrimidinyl)-;4-Amino-N-(5,6-dimethoxy-4-pyrimidinyl)benzensulfonamide;Fanasil;Ro 4-4393;Sulfadoxin;Sulformetoxin;Sulforthodimethoxine; EINECS: 219-504-9 Density: 1.441

Tetrahydrozoline hydrochloride CAS 522-48-5

Name: Tetrahydrozoline hydrochloride CAS No.:522-48-5 Formula: C13H16N2.HCl Molecular Weight : 236.74 Synonyms: 1H-Imidazole,4,5-dihydro-2-(1,2,3,4-tetrahydro-1-naphthalenyl)-, monohydrochloride (9CI);2-(1,2,3,4-Tetrahydro-1-naphthyl)-2-imidazoline hydrochloride;Murine Plus;Tinarhinin;Tyzanol hydrochloride;Tyzine

2,2,6,6-Tetramethyl-4-piperidinol CAS 2403-88-5

Name: 2,2,6,6-Tetramethyl-4-piperidinol CAS No.:2403-88-5 Formula: C9H19NO Molecular Weight : 157.25 Synonyms: NSC 16575;4-Hydroxyl-2,2,6,6-tetramethyl piperidine;DN 10;DN10 (UV absorber);HOTEMP;HTMP;Lastar A; EINECS: 219-291-2 Density: 0.891 g/cm3 Boiling Point: 214.1 °C at 760 mmHg Flash Point: 53 °C Solubility:

3-Methacryloxypropylmethyldimethoxysilane CAS 14513-34-9

Name: 3-Methacryloxypropylmethyldimethoxysilane CAS No.:14513-34-9 Formula: C10H20O4Si Molecular Weight : 232.35 Synonyms: Methacrylicacid, 3-(dimethoxymethylsilyl)propyl ester (7CI,8CI);1-Propanol,3-(dimethoxymethylsilyl)-, methacrylate (8CI);2-Propenoic acid,2-methyl-, 3-(dimethoxymethylsilyl)propyl

Spectinomycin dihydrochloride pentahydrate CAS 22189-32-8

Name: Spectinomycin dihydrochloride pentahydrate CAS No.:22189-32-8 Formula: C14H24N2O7·2(HCl)·5(H2O) Molecular Weight : 493.35 Synonyms: 4H-Pyrano[2,3-b][1,4]benzodioxin-4-one,decahydro-4a,7,9-trihydroxy-2-methyl-6,8-bis(methylamino)-, dihydrochloride,pentahydrate

2-Propenoic acid, 3-(3-(trifluoromethyl)phenyl)- CAS 779-89-5

Name: 2-Propenoic acid, 3-(3-(trifluoromethyl)phenyl)- CAS No.:779-89-5 Formula: C10H7F3O2 Molecular Weight : 216.16 Synonyms: 3-(3-(Trifluoromethyl)phenyl)-2-propenoic acid;3-(3-(Trifluoromethyl)phenyl)acrylic acid; EINECS: 212-301-6 Density: 1.363 g/cm3 Boiling Point: 289.6 °C at 760 mmHg Flash Point: 129

Ranitidine hydrochloride CAS 66357-35-5

Name: Ranitidine hydrochloride CAS No.:66357-35-5 Formula: C13H22N4O3S Molecular Weight : 314.4 Synonyms: Midaven;Zenetac;1,1-Ethenediamine, N-(2-(((5-((dimethylamino)methyl)-2-furanyl)methyl)thio)ethyl)-N'-methyl-2-nitro-;Acidex; EINECS: 266-332-5 Density: 1.184 g/cm3 Boiling Point: 437.1 °C at 760 mmHg Flash Point:

7H-Furo[3,2-g][1]benzopyran-7-one CAS 66-97-7

Name: 7H-Furo[3,2-g][1]benzopyran-7-one CAS No.:66-97-7 Formula: C11H6O3 Molecular Weight : 186.17 Synonyms: Furocoumarin(6CI);2-Propenoic acid, 3-(6-hydroxy-5-benzofuranyl)-, d-lactone;6,7-Furanocoumarin;Ficusin;Furo[2',3':7,6]coumarin;Furo[4',5':6,7]coumarin;NSC 404562;Psoralen;Psoralene; EINECS: 200-639-7 Density:

Imidazo[1,2-a]pyridine CAS 274-76-0

Name: Imidazo[1,2-a]pyridine CAS No.:274-76-0 Formula: C7H6N2 Molecular Weight : 118.14 Synonyms: 1,3a-Diazaindene;1-Azaindolizine;Pyridino[1',2':1,2]glyoxaline;Pyrimidazole; Density: 1.142 g/cm3 Boiling Point: 103oC (1 mmHg) Flash Point: 113oC

1-Propene, 3-chloro- CAS 107-05-1

Name: 1-Propene, 3-chloro- CAS No.:107-05-1 Formula: C3H5Cl Molecular Weight : 76.49 Synonyms: Propene,3-chloro- (8CI);1-Chloro-2-propene;2-Propenyl chloride;3-Chloro-1-propene;3-Chloropropene;3-Chloropropylene;NSC 20939; EINECS: 203-457-6 Density: 0.939 g/cm3 Boiling Point: 41.6 °C at 760 mmHg Flash Point:

Dutasteride CAS 164656-23-9

Name: Dutasteride CAS No.:164656-23-9 Formula: C27H30F6N2O2 Molecular Weight : 528.53 Synonyms: 4-Azaandrost-1-ene-17-carboxamide,N-[2,5-bis(trifluoromethyl)phenyl]-3-oxo-,(

Methane, trifluoro- CAS 75-46-7

Name: Methane, trifluoro- CAS No.:75-46-7 Formula: CHF3 Molecular Weight : 70.02 Synonyms: Arcton 1;Carbon trifluoride;Ecolo Ace 23;FC 23;FC 23 (fluorocarbon);FE 13;Fluoroform;Fluoryl;Freon 23;Fron 23;Genetron 23;HCFC 23;HFC 23;Methyltrifluoride;R 23;R 23 (halocarbon);Trifluoromethane (CHF3); EINECS:

Hydroxylamine CAS 7803-49-8

Name: Hydroxylamine CAS No.:7803-49-8 Formula: H3N O Molecular Weight : 33.04 Synonyms: HDA;Hydroxyamine; Telclean 1200A EINECS: 232-259-2 Density: 1.078 Boiling Point: 56.5 °C at 760 mmHg Flash Point: °C Solubility: soluble in cold water, decomposes in hot water Appearance: white needles or flakes

Glycyrrhizic acid CAS 1405-86-3

Name: Glycyrrhizic acid CAS No.:1405-86-3 Formula: C42H62O16 Molecular Weight : 823.04 Synonyms: 6-[(11-carboxy-4,4,6a,6b,8a,11,14b-heptamethyl-14-oxo-2,3,4a,5,6,7,8,9,10,12,12a,14a-dodecahydro-1H-picen-3-yl)oxy]-5-(6-carboxy-3,4,5-trihydroxy-oxan-2-yl)oxy-3,4-dihydroxy-oxane-2-carboxylic acid;See

Letrozole CAS 112809-51-5

Name: Letrozole CAS No.:112809-51-5 Formula: C17H11N5 Molecular Weight : 285.30 Synonyms: 4,4'-(1H-1,2,4-Triazol-1-ylmethylene)bisbenzonitrile Density: 1.21g/cm3 Boiling Point: 374.4 °C at 760 mmHg Flash Point: 180.3 °C Appearance: white to off-whitecrystalline powder

Pigment Violet 3 CAS 1325-82-2

Name: Pigment Violet 3 CAS No.:1325-82-2 Formula: C20H19N3 Molecular Weight : 301.385 Synonyms: C.I.Pigment Violet 3 (8CI);Benzoic Methyl Violet Lake;Chemisperse CV 8030;Conc. Violet R;Fanal Violet R Supra;Fanal Violet RA Supra;Fast Bronze Violet;Fast Violet Toner R;Fastel Violet R;Fastel Violet R Supra;Halopont

Lansoprazole CAS 103577-45-3

Name: Lansoprazole CAS No.:103577-45-3 Formula: C16H14F3N3O2S Molecular Weight : 369.36 Synonyms: Amarin;Promp;Monolitum;Lanz;Prevacid (TN);Ulpax;Lansoprazolum [INN-Latin];Takepron;Blason;2-(((3-Methyl-4-(2,2,2-trifluoroethoxy)-2-pyridyl)methyl)sulfinyl)benzimidazole;Ilsatec;A 65006;Bamalite;Prevacid

Benzenesulfonic acid,[[4-[bis[4-[(sulfophenyl)amino]phenyl]methylene]-2,5-cyclohexadien-1-ylidene]amino]-,sodium salt (1:2) CAS 28983-56-4

Name: Benzenesulfonic acid,[[4-[bis[4-[(sulfophenyl)amino]phenyl]methylene]-2,5-cyclohexadien-1-ylidene]amino]-,sodium salt (1:2) CAS No.:28983-56-4 Formula: C37H27N3Na2O9S3 Molecular Weight : 799.79 Synonyms: Benzenesulfonicacid,

Benzene,1-bromo-3-(methylthio)- CAS 33733-73-2

Name: Benzene,1-bromo-3-(methylthio)- CAS No.:33733-73-2 Formula: C7H7BrS Molecular Weight : 203.1 Synonyms: Sulfide,m-bromophenyl methyl (6CI,7CI,8CI);1-Bromo-3-(methylsulfanyl)benzene;1-Bromo-3-(methylthio)benzene;3-(Methylthio)bromobenzene;3-Bromophenyl methylsulfide;3-Bromothioanisole;Methyl m-bromophenyl

Acetyl-L-Carnitine CAS 14992-62-2

Name: 1-Propanaminium,2-(acetyloxy)-3-carboxy-N,N,N-trimethyl-, inner salt CAS No.:14992-62-2 Formula: C9H17NO4 Molecular Weight : 203.24 Synonyms: Ammonium,(57263500,3-carboxy-2-hydroxypropyl)trimethyl-, hydroxide, inner salt, acetate (8CI);Vitamin BT, acetate (6CI);Acetylcarnitine;Carnitineacetyl

Cephalexin CAS 15686-71-2

Name: Cephalexin CAS No.:15686-71-2 Formula: C16H17N3O4S Molecular Weight : 347.39 Synonyms: Sencephalin;Septilisin;Sialexin;Sinthecillin;Sporidex;Syncle;Tepaxin;Uphalexin;Vetolexin;Winlex;Alcephin;Biocef ;Cefablan;Cefadina;Cefalexin;Cefaloto;Ceforal;Celexin;Cepexin; EINECS: 239-773-6 Density: 1.5 g/cm3 Boiling Point:

5-Bromo-2-pyridinecarbonitrile CAS 97483-77-7

Name: 5-Bromo-2-pyridinecarbonitrile CAS No.:97483-77-7 Formula: C6H3BrN2 Molecular Weight : 183.01 Synonyms: 2-Cyano-5-Bromopyridin;2-Cyano-5-Bromo pyridine;5-Bromo-2-cyanopyridine 97483-77-7;LDP05:5-Bromo-2-Cyanopyridine;5-bromopicolinonitrile;5-bromopyridine-2-carbonitrile;5-Bromo-2-cyanopyridine; EINECS:

Calcium phosphate CAS 7758-87-4

Name: Calcium phosphate CAS No.:7758-87-4 Formula: Ca3O8P2 Molecular Weight : 310.18 Synonyms: Posture (body position)Posture (calcium supplement);Bonarka;Tricalcium bis(orthophosphate), with a fluorine content of less than 0,005 % by weight on the dry anhydrous product;Calcigenol simple;Calcium orthophosphate,

L(-)-Carnitine hydrochloride CAS 6645-46-1

Name: L(-)-Carnitine hydrochloride CAS No.:6645-46-1 Formula: C7H16ClNO3 Molecular Weight : 196.68 Synonyms: 1-Propanaminium,3-carboxy-2-hydroxy-N,N,N-trimethyl-, chloride,(57263492,R)-;Ammonium,(57263493,3-carboxy-2-hydroxypropyl)trimethyl-,chloride,(57263494,-)- (8CI);(R)-(-)-Carnitine

Methanone,[6-hydroxy-2-(4-hydroxyphenyl)benzo[b]thien-3-yl][4-[2-(1-piperidinyl)ethoxy]phenyl]- CAS 84449-90-1

Name: Methanone,[6-hydroxy-2-(4-hydroxyphenyl)benzo[b]thien-3-yl][4-[2-(1-piperidinyl)ethoxy]phenyl]- CAS No.:84449-90-1 Formula: C28H27NO4S Molecular Weight : 473.58 Synonyms: Keoxifene;LY 139481;[2-(4-Hydroxyphenyl)-6-hydroxybenzo[b]thien-3-yl][4-(2-(1-piperidinyl)ethoxy)phenyl]methanone; Density: 1.289

C.I. Acid Brown 282 CAS 12219-65-7

Name: C.I. Acid Brown 282 CAS No.:12219-65-7 Formula: Unspecified Molecular Weight : 0 Synonyms: Acid Brown S-GR;DyalanBrown S-GR;Everacid Brown GR;Everlan Brown GR;IrgadermBrown M;Irgaderm Brown M 210;Irgasperse Brown 4R-U;Irgasperse Brown 4RL-U;Lanacron Brown S-GR;Lanacron DarkBrown S-GR;Lanasyn Brown S-Bgl;Levaderm

Pendimethalin CAS 40487-42-1

Name: Pendimethalin CAS No.:40487-42-1 Formula: C13H19N3O4 Molecular Weight : 281.35 Synonyms: 3,4-Dimethyl-2,6-dinitro-N-(1-ethylpropyl)aniline;3,4-Xylidine, 2,6-dinitro-N-(1-ethylpropyl)-;Benzenamine, 3,4-dimethyl-2,6-dinitro-N-(1-ethylpropyl)-;N-(1-Aethylpropyl)-2,6-dinitro-3,4-xylidin;Phenoxalin; EINECS:

Benzenamine,2-bromo- CAS 615-36-1

Name: Benzenamine,2-bromo- CAS No.:615-36-1 Formula: C6H6BrN Molecular Weight : 172.03 . Synonyms: Aniline,o-bromo- (8CI);2-Amino-1-bromobenzene;2-Bromobenzenamine;2-Bromophenylamine;NSC 7086;o-Aminobromobenzene;o-Bromoaniline; EINECS: 210-421-3 Density: 1.594 g/cm3 Boiling Point: 227 °C at 760 mmHg Flash Point: 91.1

4-Bromoaniline CAS 106-40-1

Name: 4-Bromoaniline CAS No.:106-40-1 Formula: C6H6BrN Molecular Weight : 172.04 Synonyms: Aniline,p-bromo- (8CI);1-Amino-4-bromobenzene;benzenamine, 4-bromo-;Aniline, p-bromo-;p-Bromophenylamine;p-Bromoaniline; EINECS: 203-393-9 Density: 1.594 g/cm3 Boiling Point: 225.9 °C at 760 mmHg Flash Point: 90.4

3,5-Dibromopyridine CAS 625-92-3

Name: 3,5-Dibromopyridine CAS No.:625-92-3 Formula: C5H3Br2N Molecular Weight : 236.89 Synonyms: Pyridine, 3,5-dibromo-; EINECS: 210-916-4 Density: 2.059 g/cm3 Boiling Point: 222 °C at 760 mmHg Flash Point: 84.3 °C Appearance: Light-Yellow Solid

1-Bromo-2,5-difluorobenzene CAS 399-94-0

Name: 1-Bromo-2,5-difluorobenzene CAS No.:399-94-0 Formula: C6H3BrF2 Molecular Weight : 192.99 Synonyms: 2,5-Difluoro-1-bromobenzene;2,5-Difluorobromobenzene;2-Bromo-1,4-difluorobenzene;NSC 10250; EINECS: 206-920-0 Density: 1.692 g/cm3 Boiling Point: 155.4 °C at 760 mmHg Flash Point: 65 °C Solubility: insoluble in

2-Pyridinecarbonitrile, 4-chloro- CAS 19235-89-3

Name: 2-Pyridinecarbonitrile, 4-chloro- CAS No.:19235-89-3 Formula: C6H3ClN2 Molecular Weight : 138.55 Synonyms: Picolinonitrile,4-chloro- (6CI,8CI);2-Cyano-4-chloropyridine;4-Chloro-2-cyanopyridine;4-Chloro-2-pyridinecarbonitrile;4-Chloro-pyridine-2-carbonitrile;4-Chloropicolinonitrile; EINECS: -0 Density: 1.3±0.1

3-Butenenitrile CAS 109-75-1

Name: 3-Butenenitrile CAS No.:109-75-1 Formula: C4H5N Molecular Weight : 67.10 Synonyms: 1-Butene-3-nitrile;1-Propene-3-nitrile;2-Propenyl cyanide;3-Butenonitrile;3-Butenylnitrile;3-Cyano-1-propene;3-Cyanopropene;Allyl cyanide;NSC 2583;Vinylacetonitrile;b-Butenonitrile; EINECS: 203-701-1 Density: 0.807 g/cm3 Boiling

Magnesium fluosilicate CAS 16949-65-8

Name: Magnesium fluosilicate CAS No.:16949-65-8 Formula: MgSiF6 Molecular Weight : 166.38 Synonyms: Magnesiumfluosilicate (6CI);Silicate(2-), hexafluoro-, magnesium (8CI);Fluosilicicacid magnesium salt;Magnesium hexafluorosilicate;Magnesium hexafluorosilicate(2-);Magnesium silicofluoride;Silicon fluoridemagnesium

2-Cyaniminothiazolidine CAS 26364-65-8

Name: 2-Cyaniminothiazolidine CAS No.:26364-65-8 Formula: C4H5N3S Molecular Weight : 127.16 Synonyms: Cyanamide,(57263479,4,5-dihydro-2-thiazolyl)- (9CI);D2,N-Thiazolidinecarbamonitrile (8CI);2-(Cyanoimino)thiazolidine;2-Cyanoimino-1,3-thiazolidine;2-Thiazolidinylidenecyanamide; EINECS: 427-720-1 Density: 1.43

3-Fluorobenzotrifluoride CAS 401-80-9

Name: 3-Fluorobenzotrifluoride CAS No.:401-80-9 Formula: C7H4F4 Molecular Weight : 164.10 Synonyms: Toluene,m,a,a,a-tetrafluoro- (6CI,7CI,8CI);1-Fluoro-3-(trifluoromethyl)benzene;3-Fluoro(trifluoromethyl)benzene;Benzene,1-fluoro-3-(trifluoromethyl)-;NSC

Phenol,2,5-dimethyl- CAS 95-87-4

Name: Phenol,2,5-dimethyl- CAS No.:95-87-4 Formula: C8H10O Molecular Weight : 122.18 Synonyms: 2,5-Xylenol(8CI);1,2,5-Xylenol;1-Hydroxy-2,5-dimethylbenzene;2,5-Dimethylphenol;2-Hydroxy-p-xylene;3,6-Dimethylphenol;NSC 2599;p-Xylenol; EINECS: 202-461-5 Density: 1.041 g/cm3 Boiling Point: 211.1 °C at 760 mmHg Flash

Bicyclo[2.2.1]heptane-7-methanesulfonicacid, 2-bromo-4,7-dimethyl-3-oxo-, ammonium salt (1:1),(57263472,1S,2S,4R,7R)- CAS 14575-84-9

Name: Bicyclo[2.2.1]heptane-7-methanesulfonicacid, 2-bromo-4,7-dimethyl-3-oxo-, ammonium salt (1:1),(57263473,1S,2S,4R,7R)- CAS No.:14575-84-9 Formula: C10H18BrNO4S Molecular Weight : 328.22 Synonyms: 8-Bornanesulfonicacid, 3-bromo-2-oxo-, ammonium salt,(57263474,+)- (8CI);Bicyclo[2.2.1]heptane-7-methanesulfonic

Vancomycin CAS 1404-90-6

Name: Vancomycin CAS No.:1404-90-6 Formula: C66H75Cl2N9O24 Molecular Weight : 1449.40 Synonyms: 22H-8,11:18,21-Dietheno-23,36-(iminomethano)-13,16:31,35-dimetheno-1H,16H-[1,6,9]oxadiazacyclohexadecino[4,5-m][10,2,16]benzoxadiazacyclotetracosine-26-carboxylicacid,

Acetic acid,2-[2-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]ethoxy]-, hydrochloride(1:2) CAS 83881-52-1

Name: Acetic acid,2-[2-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]ethoxy]-, hydrochloride(1:2) CAS No.:83881-52-1 Formula: C21H25ClN2O3.2(HCl) Molecular Weight : 461.81 Synonyms: Alercet;Alerid;Alerlisin;Cesta;Cetirizinedihydrochloride;Cetirizine

2-Mercaptomethylbenzimidazole CAS 53988-10-6

Name: 2-Mercaptomethylbenzimidazole CAS No.:53988-10-6 Formula: C8H8N2S Molecular Weight : 164.24 Synonyms: 4(or 5)-Methyl-2-benzimidazolethiol; EINECS: 258-904-8 Density: 1.25 Boiling Point: 276.8oC at 760mmHg Flash Point: 121.2oC

1,3,5-Triazine-2,4-diamine,6-chloro-N2,N4-bis(1-methylethyl)- CAS 139-40-2

Name: 1,3,5-Triazine-2,4-diamine,6-chloro-N2,N4-bis(1-methylethyl)- CAS No.:139-40-2 Formula: C9H16 Cl N5 Molecular Weight : 229.7098 Synonyms: 1,3,5-Triazine-2,4-diamine,6-chloro-N,N'-bis(1-methylethyl)- (9CI); s-Triazine, 2-chloro-4,6-bis(isopropylamino)-(6CI,8CI);

1,1-Ethenediamine,N-[(6-chloro-3-pyridinyl)methyl]-N-ethyl-N'-methyl-2-nitro- CAS 120738-89-8

Name: 1,1-Ethenediamine,N-[(6-chloro-3-pyridinyl)methyl]-N-ethyl-N'-methyl-2-nitro- CAS No.:120738-89-8 Formula: C11H15ClN4O2 Molecular Weight : 270.72 Synonyms: N-[(6-chloropyridin-3-yl)methyl]-N-ethyl-N'-methyl-2-nitroethene-1,1-diamine;Capstar; Density: 1.254 g/cm3 Boiling Point: 417.2 °C at 760 mmHg Flash Point:

1-Naphthalenesulfonicacid, 8-(phenylamino)- CAS 82-76-8

Name: 1-Naphthalenesulfonicacid, 8-(phenylamino)- CAS No.:82-76-8 Formula: C16H13NO3S Molecular Weight : 299.34 Synonyms: 1-Naphthalenesulfonicacid, 8-anilino- (7CI,8CI);1-(Phenylamino)-8-naphthalenesulfonic acid;1-Anilino-8-naphthalenesulfonic

3-Pyridinecarboxylicacid,(57263463,2R)-3,4-dihydro-2,5,7,8-tetramethyl-2-[(4R,8R)-4,8,12-trimethyltridecyl]-2H-1-benzopyran-6-ylester, rel- CAS 51898-34-1

Name: 3-Pyridinecarboxylicacid,(57263464,2R)-3,4-dihydro-2,5,7,8-tetramethyl-2-[(4R,8R)-4,8,12-trimethyltridecyl]-2H-1-benzopyran-6-ylester, rel- CAS No.:51898-34-1 Formula: C35H53NO3 Molecular Weight :

3-Pyridinol, 6-amino- CAS 55717-46-9

Name: 3-Pyridinol, 6-amino- CAS No.:55717-46-9 Formula: C5H6N2O Molecular Weight : 110.11 Synonyms: 2-Amino-5-pyridinol;6-Amino-3-hydroxypyridine;6-Amino-3-pyridinol; EINECS: -0 Density: 1.32 g/cm3 Boiling Point: 447.4 °C at 760 mmHg Flash Point: 224.4 °C Appearance: light yellow crystal

4-Methylthiazole CAS 693-95-8

Name: 4-Methylthiazole CAS No.:693-95-8 Formula: C4H5NS Molecular Weight : 99.15 Synonyms: 4-Methyl-1,3-thiazole;NSC 42976;Thiazole, 4-methyl-; EINECS: 211-764-1 Density: 1.121 g/cm3 Boiling Point: 134.3 °C at 760 mmHg Flash Point: 37.3 °C Appearance: Clear colourless to slightly yellow liquid

2-Amino-5-bromo-4-methyl-3-nitropyridine CAS 100367-40-6

Name: 2-Amino-5-bromo-4-methyl-3-nitropyridine CAS No.:100367-40-6 Formula: C6H6BrN3O2 Molecular Weight : 233.0421 Synonyms: 4-Picoline,2-amino-5-bromo-3-nitro- (6CI);2-Amino-5-bromo-3-nitro-4-methylpyridine;5-Bromo-4-methyl-3-nitropyridin-2-amine; EINECS: -0 Density: 1.795 g/cm3 Boiling Point: 314.1 °C at 760

5H-Dibenz[b,f]azepine-5-propanamine,3-chloro-10,11-dihydro-N,N-dimethyl-, hydrochloride (1:1) CAS 17321-77-6

Name: 5H-Dibenz[b,f]azepine-5-propanamine,3-chloro-10,11-dihydro-N,N-dimethyl-, hydrochloride (1:1) CAS No.:17321-77-6 Formula: C19H23ClN2.HCl Molecular Weight : 351.32 Synonyms: 5H-Dibenz[b,f]azepine,3-chloro-5-[3-(dimethylamino)propyl]-10,11-dihydro-, monohydrochloride (8CI);5H-Dibenz[b,f]azepine-5-propanamine,

Misoprostol CAS 59122-46-2

Name: Misoprostol CAS No.:59122-46-2 Formula: C22H38O5 Molecular Weight : 382.60 Synonyms: Cytotec;Isprelor;Misogon;Misoprost;Misoprostil;SC 29333; Density: 1.078 g/cm3 Boiling Point: 497.3 °C at 760 mmHg Flash Point: 160.3 °C Appearance: water-soluble, viscous liquid

6-Aminopenicillanic acid CAS 551-16-6

Name: 6-Aminopenicillanic acid CAS No.:551-16-6 Formula: C8H12N2O3S Molecular Weight : 216.25 Synonyms: 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylicacid, 6-amino-3,3-dimethyl-7-oxo- (6CI,8CI);4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid,6-amino-3,3-dimethyl-7-oxo-, [2S-(2a,5a,6b)]-;(+)-6-Aminopenicillanic

Naphthalene,1-fluoro- CAS 321-38-0

Name: Naphthalene,1-fluoro- CAS No.:321-38-0 Formula: C10H7F Molecular Weight : 146.17 Synonyms: 1-Fluoronaphthalene;NSC 4690;a-Fluoronaphthalene; EINECS: 206-287-0 Density: 1.144 g/cm3 Boiling Point: 215 °C at 760 mmHg Flash Point: 65.6 °C Appearance: clear slightly yellow to yellow-brown liquid

Argon CAS 7440-37-1

Name: Argon CAS No.:7440-37-1 Formula: Ar Molecular Weight : 39.95 Synonyms: 40Ar;Argon, isotope of mass 40; Argon-40; Khladon R 740; R 740 EINECS: 231-147-0 Density: 1.784（0ºC） Boiling Point: -186 ºC Solubility: slight

4-Bromo-2-fluorobenzyl bromide CAS 76283-09-5

Name: 4-Bromo-2-fluorobenzyl bromide CAS No.:76283-09-5 Formula: C7H5Br2F Molecular Weight : 267.92 Synonyms: 4-Bromo-1-(bromomethyl)-2-fluorobenzene;4-bromo-1-(bromomethyl)-2-fluoro-benzene;4-Bromo 2-fluorobenzylbromide;2-Fluoro-bromobenzyl bromide;4-Bromo-2-fluorobenzylbromide;3,4-difluoro benzyl

DL-Mandelic acid CAS 611-72-3

Name: DL-Mandelic acid CAS No.:611-72-3 Formula: C8H8O3 Molecular Weight : 152.15 Synonyms: DL-2-Hydroxy-2-phenylacetic acid;(+/-)-alpha-Hydroxyphenylacetic acid; EINECS: 210-277-1 Density: 1.322 g/cm3 Boiling Point: 321.827 °C at 760 mmHg Flash Point: 162.649 °C Solubility: 150 g/L (20℃)

Phosphate(1-),hexafluoro-, lithium (1:1) CAS 21324-40-3

Name: Phosphate(1-),hexafluoro-, lithium (1:1) CAS No.:21324-40-3 Formula: F6LiP Molecular Weight : 151.90 Synonyms: Lithiumhexafluorophosphate (7CI);Phosphate(1-), hexafluoro-, lithium (8CI,9CI);Lithium fluophosphate;Lithium hexafluorophosphate (LiPF6);Lithiumhexafluorophosphate(1-);Lithium phosphorus fluoride

Emamectin, benzoate (salt) CAS 137512-74-4

Name: Emamectin, benzoate (salt) CAS No.:137512-74-4 Formula: C48H73NO13.C7H6O2 Molecular Weight : 1008.24000 Synonyms: (4"R)-4"-Deoxy-4"-(methylamino)-avermectin B1 benzoate;Methylamino abamectin benzoate(salt);

1,4-Benzenediamine,N1-(4-aminophenyl)- CAS 537-65-5

Name: 1,4-Benzenediamine,N1-(4-aminophenyl)- CAS No.:537-65-5 Formula: C12H13 N3 Molecular Weight : 199.25 Synonyms: 1,4-Benzenediamine,N-(4-aminophenyl)- (9CI); Diphenylamine, 4,4'-diamino- (6CI,7CI,8CI);4,4'-Diaminodiphenylamine; 4,4'-Iminodianiline; Benzenamine, 4,4'-iminobis-;Bis(4-aminophenyl)amine;

Benzenamine,N-ethyl-N-[2-[1-(2-methylpropoxy)ethoxy]ethyl]-4-(2-phenyldiazenyl)- CAS 34432-92-3

Name: Benzenamine,N-ethyl-N-[2-[1-(2-methylpropoxy)ethoxy]ethyl]-4-(2-phenyldiazenyl)- CAS No.:34432-92-3 Formula: C22H31 N3 O2 Molecular Weight : 369.50044 Synonyms: Acetaldehyde,isobutyl 2-[N-ethyl-p-(phenylazo)anilino]ethyl acetal (8CI); Benzenamine,N-ethyl-N-[2-[1-(2-methylpropoxy)ethoxy]ethyl]-4-(phenylazo)-

Reboxetine mesylate CAS 98769-84-7

Name: Reboxetine mesylate CAS No.:98769-84-7 Formula: C19H23NO3.CH4O3S Molecular Weight : 409.50 Synonyms: Reboxetine mesylate [USAN];(2R)-2-[(R)-(2-ethoxyphenoxy)-phenyl-methyl]morpholine; methanesulfonic acid;FCE 20124;Morpholine, 2-((R)-(2-ethoxyphenoxy)phenylmethyl)-,(57263446,2R)-rel-,

Escitalopram oxalate CAS 219861-08-2

Name: Escitalopram oxalate CAS No.:219861-08-2 Formula: C22H23FN2O5 Molecular Weight : 414.43 Synonyms: Escitapram;S-(+)-1-[3-(dimethyl-amino)propyl]-1-(p-fluorophenyl)-5-phthalancarbonitrile oxalate;S-Citalopram Oxalate(escitalopram oxalate);(S)-Citalopram oxalate;5-Isobenzofurancarbonitrile,

Narirutin CAS 14259-46-2

Name: 4H-1-Benzopyran-4-one,7-[[6-O-(6-deoxy-a-L-mannopyranosyl)-b-D-glucopyranosyl]oxy]-2,3-dihydro-5-hydroxy-2-(4-hydroxyphenyl)-,(57263441,2S)- CAS No.:14259-46-2 Formula: C27H32O14 Molecular Weight :

Vitamin d CAS 511-28-4

Name: VitaminD4 CAS No.:511-28-4 Formula: C28H46O Molecular Weight : 398.66 Synonyms: 9,10-Secoergosta-5,7,10(19)-trien-3-ol,(57263439,3b,5Z,7E)- (9CI);Ergocalciferol,22,23-dihydro- (6CI);Vitamin D4 (7CI,8CI);22,23-Dihydroercalciol;22,23-Dihydroergocalciferol;22,23-Dihydrovitamin

D-(+)-Methyl-alpha-(2-thienylethamino)(2-chlorophenyl)acetate hydrochloride CAS 141109-19-5

Name: D-(+)-Methyl-alpha-(2-thienylethamino)(2-chlorophenyl)acetate hydrochloride CAS No.:141109-19-5 Formula: C15H17Cl2NO2S Molecular Weight : 346.272 Synonyms: Benzeneaceticacid, 2-chloro-a-[[2-(2-thienyl)ethyl]amino]-,methyl ester, hydrochloride,(57263436,S)-;(+)-(S)-Methyl

1-Benzyl-5-phenylbarbituric acid CAS 72846-00-5

Name: 1-Benzyl-5-phenylbarbituric acid CAS No.:72846-00-5 Formula: C17H14N2O3 Molecular Weight : 294.30 Synonyms: 5-Phenyl-1-benzylbarbituric acid; EINECS: 276-940-2 Density: 1.304 g/cm3 Solubility: insoluble in water Appearance: Odorless powder

Benzamide, 3-chloro- CAS 618-48-4

Name: Benzamide, 3-chloro- CAS No.:618-48-4 Formula: C7H6ClNO Molecular Weight : 155.58 Synonyms: Benzamide,m-chloro- (6CI,7CI,8CI);3-Chlorobenzamide;m-Chlorobenzamide; EINECS: 210-554-7 Density: 1.295 g/cm3 Boiling Point: 273.7 °C at 760 mmHg Flash Point: 119.4 °C Appearance: white powder

5H-Dibenz[b,f]azepine-5-propanamine,3-chloro-10,11-dihydro-N,N-dimethyl-, hydrochloride (1:1) CAS 303-49-1

Name: 5H-Dibenz[b,f]azepine-5-propanamine,3-chloro-10,11-dihydro-N,N-dimethyl- CAS No.:303-49-1 Formula: C19H23ClN2 Molecular Weight : 314.89 Synonyms: 5H-Dibenz[b,f]azepine,3-chloro-5-[3-(dimethylamino)propyl]-10,11-dihydro-

Thiabendazole CAS 148-79-8

Name: Thiabendazole CAS No.:148-79-8 Formula: C10H7N3S Molecular Weight : 201.25 Synonyms: Thibenzole 200;Mertect 160;MK 360;Tiabenda;Benzimidazole, 2- (4-thiazolyl)-;E-Z-Ex;4-(2-benzimidazolyl)thiazole;Polival;Hokustar

Chloroacetonitrile CAS 107-14-2

Name: Chloroacetonitrile CAS No.:107-14-2 Formula: C2H2ClN Molecular Weight : 75.50 Synonyms: Acetonitrile,chloro- (6CI,8CI,9CI);2-Chloroacetonitrile;Chloromethyl cyanide;Cyanomethyl chloride;Monochloroacetonitrile;Acetonitrile,2-chloro-;a-Chloroacetonitrile; EINECS: 203-467-0 Density: 1.138 g/cm3 Boiling Point: 126.5

Poly(acrylamide-co-diallyldimethylammonium chloride) CAS 26590-05-6

Name: Poly(acrylamide-co-diallyldimethylammonium chloride) CAS No.:26590-05-6 Formula: (C8H16ClN)n.(C3H5NO)m Molecular Weight : 232.75000 Synonyms: Agequat C 3204;Merquat 2200;CV 5380;XB 54-15-1;WT 2640;Polyquaternium-7,26590-05-6;Himacs SC 100;Polyquaternium 7;WT 5504;Hydraid 777;PAS-J 11;Merquat S;Merquat 550;E

Ethisterone CAS 434-03-7

Name: Ethisterone CAS No.:434-03-7 Formula: C21H28O2 Molecular Weight : 312.49 Synonyms: Gestoral;Testosterone, 17-ethynyl-;17-Hydroxy-17-alpha-pregn-4-en-20-yn-3-one;Trosinone;Ora-Lutin;Etisterona [INN-Spanish];Progestin P;Lutocylol;17-alpha-Pregn-4-en-20-yn-3-one, 17-hydroxy-;Ethinyltestosterone;Ethinyl

C.I. Vat Blue 43 CAS 1327-79-3

Name: C.I. Vat Blue 43 CAS No.:1327-79-3 Formula: C18H12N2O Molecular Weight : 274.32 Synonyms: Hydron Blue 3R Stabilosol;Hydron Blue 3RC;Hydron Blue R;Sulfurized Vat Blue R-T;4-(9H-Carbazol-3-ylimino)cyclohexa-2,5-dien-1-one; EINECS: 215-499-2 Density: 1.26 g/cm3 Boiling Point: 494.643 °C at 760 mmHg Flash Point:

Thiosalicylic acid CAS 147-93-3

Name: Thiosalicylic acid CAS No.:147-93-3 Formula: C7H6O2S Molecular Weight : 154.18 . Synonyms: USAF XR-35;Benzoic acid, 2-mercapto-, monosodium salt;o-Mercaptobenzoesaeure [German];Salicylic acid, 2-thio-;(6-sulfanylidene-1-cyclohexa-2,4-dienylidene)methanediolate;o-Sulfhydrylbenzoic acid;Benzoic acid,

9-Fluorenone CAS 486-25-9

Name: 9-Fluorenone CAS No.:486-25-9 Formula: C13H8O Molecular Weight : 180.21 Synonyms: Fluoren-9-one(8CI);Fluorenone;NSC 5181; EINECS: 207-630-7 Density: 1.244 g/cm3 Boiling Point: 341.499 °C at 760

mmHg Flash Point: 144.142 °C Solubility: insoluble in water Appearance: yellow flakes, chips or crystalline powder

2,3-Butanediol,(57263419,2R,3R)- CAS 24347-58-8

Name: 2,3-Butanediol,(57263420,2R,3R)- CAS No.:24347-58-8 Formula: C4H10O2 Molecular Weight : 90.12 Synonyms: 2,3-Butanediol, [R-(R*,R*)]-

1-Adamantanol CAS 768-95-6

Name: 1-Adamantanol CAS No.:768-95-6 Formula: C10H16O Molecular Weight : 152.24 Synonyms: Tricyclo(3.3.1.1(sup 3,7))decan-1-ol (9CI);1-Adamantol;Tricyclo[3.3.1.13,7]decan-1-ol;Adapalene-1;1-Ad-OH;Tricyclo(3.3.1.13,7)decan-1-ol;1-Hydroxyadamantane;adamantan-1-ol;tricyclo[3.3.1.1~3,7~]decan-1-ol; EINECS:

Enduracidin CAS 11115-82-5

Name: Enduracidin CAS No.:11115-82-5 Formula: Unspecified Molecular Weight : 2340.2677 Synonyms:Enramycin;Enradin;

Pravastatin sodium CAS 81131-70-6

Name: Pravastatin sodium CAS No.:81131-70-6 Formula: C23H35NaO7 Molecular Weight : 334.57 Synonyms: Mevalotin;Pravachol;Pravachol (TN);Pravastatine [French];1-Naphthaleneheptanoic acid,1,2,6,7,8,8a-hexahydro-a,?,6-trihydroxy- 2-methyl-8-[(2S)-2-methyl-1-oxobutoxy]-,monosodium salt,(

Sodium tripolyphosphate CAS 7758-29-4

Name: Sodium tripolyphosphate CAS No.:7758-29-4 Formula: Na5P3O10 Molecular Weight : 367.86 Synonyms: 7758-29-4 EINECS: 231-838-7 Density: >1.5 g/cm3 (20oC) Solubility: 20 g/100 mL (20°C) in water Appearance: white or colourless crystals, granules or powder

4-Nitrophthalonitrile CAS 31643-49-9

Name: 4-Nitrophthalonitrile CAS No.:31643-49-9 Formula: C8H3N3O2 Molecular Weight : 173.13 Synonyms: Phthalonitrile,4-nitro- (6CI,8CI);1,2-Dicyano-4-nitrobenzene;3,4-Dicyanonitrobenzene;4-Nitro-1,2-benzenedicarbonitrile;4-Nitro-1,2-dicyanobenzene;4-Nitro-1,2-phthalonitrile;4-Nitro-o-benzenedinitrile; EINECS:

C.I. Acid Black 107 CAS 12218-96-1

Name: C.I. Acid Black 107 CAS No.:12218-96-1 Formula: Unspecified Molecular Weight : 0 Synonyms: Acid Black 158;Acid Black BGL;Aluminium Black LBB;Anadurm Black S-BGL;Black LLB for Aluminium;C.I. Acid Black 158;Capracyl Black BGL;Cibalan BlackBGL;Covalene Black BGL;Dinalan Black BGL;Dyalan Black BGL;Fabrilan

16-Dehydropregnenolone acetate CAS 979-02-2

Name: 16-Dehydropregnenolone acetate CAS No.:979-02-2 Formula: C23H32O3 Molecular Weight : 356.50 Synonyms: 20-oxopregna-5,16-dien-3-β-yl acetate;16-Denyprasterone Acetate;16-Denyprasterone Acetate(16-DPA);16-Dehydropregnenolone Acetate(16-DPA);16 Dehydropregneninolone, Acetate;Pregna-5,16-dien-20-one, 3-

4-Pyridinecarboxylicacid, 3-bromo- CAS 13959-02-9

Name: 4-Pyridinecarboxylicacid, 3-bromo- CAS No.:13959-02-9 Formula: C6H4BrNO2 Molecular Weight : 202.00 Synonyms: Isonicotinicacid, 3-bromo- (7CI,8CI);3-Bromo-4-pyridinecarboxylic acid; EINECS: -0 Density: 1.813 g/cm3 Boiling Point: 403.1 at 760 mmHg Flash Point: 197.6 °C

Palladium oxide (PdO) CAS 1314-08-5

Name: Palladium oxide (PdO) CAS No.:1314-08-5 Formula: OPd Molecular Weight : 122.42 Synonyms: Palladium(II) oxide; EINECS: 215-218-3 Density: 8.7 g/mL at 25 °C(lit.) Solubility: practically insoluble in water Appearance: grey to black powder

9,12-Octadecadienoicacid (9Z,12Z)- CAS 60-33-3

Name: 9,12-Octadecadienoicacid (9Z,12Z)- CAS No.:60-33-3 Formula: C18H32O2 Molecular Weight : 280.50 Synonyms: 9,12-Octadecadienoicacid (Z,Z)-;Linoleic acid (8CI);(9Z,12Z)-9,12-Octadecadienoic acid;(Z,Z)-9,12-Octadecadienoic acid;9,12-Octadecadienoic acid,(57263402,Z,Z)-;9-cis,12-cis-Linoleic acid;9Z,12Z-Linoleic

11H-Dibenzo[b,e][1,4]diazepin-11-one,8-chloro-5,10-dihydro- CAS 50892-62-1

Name: 11H-Dibenzo[b,e][1,4]diazepin-11-one,8-chloro-5,10-dihydro- CAS No.:50892-62-1 Formula: C13H9ClN2O Molecular Weight : 244.68 Synonyms: 8-Chloro-5H-dibenzo[b,e][1,4]diazepin-11(10H)-one; Density: 1.327 g/cm3 Boiling Point: 334.4 °C at 760 mmHg Flash Point: 156 °C Appearance: white solid

2-Amino-4-bromoacetophenone hydrochloride CAS 5467-72-1

Name: 2-Amino-4-bromoacetophenone hydrochloride CAS No.:5467-72-1 Formula: C8H9BrClNO Molecular Weight : 250.52 Synonyms: Acetophenone,2-amino-4'-bromo-, hydrochloride (6CI,7CI,8CI);Ethanone,2-amino-1-(4-bromophenyl)-, hydrochloride (9CI);2-Amino-1-(4-bromophenyl)ethan-1-one

D-Fructose,1,6-bis(dihydrogen phosphate) CAS 488-69-7

Name: D-Fructose,1,6-bis(dihydrogen phosphate) CAS No.:488-69-7 Formula: C6H14O12P2 Molecular Weight : 340.12 Synonyms: Fructose,1,6-bis(dihydrogen phosphate), D- (6CI,8CI);D-Fructose 1,6-biphosphate;D-Fructose 1,6-bisphosphate;D-Fructose 1,6-diphosphate;Diphosphofructose;Esafosfan;Esafosfina;FDP;Fosfructose;Fructose

3,5-Dimethoxybenzaldehyde CAS 7311-34-4

Name: 3,5-Dimethoxybenzaldehyde CAS No.:7311-34-4 Formula: C9H10O3 Molecular Weight : 166.18 Synonyms: Benzaldehyde, 3,5-dimethoxy-; EINECS: 230-772-6 Density: 1.114 g/cm3 Boiling Point: 276.5 °C at 760 mmHg Flash Point: 119 °C Solubility: insoluble in water Appearance: white to beige crystalline solid

Naringenin CAS 480-41-1

Name: Naringenin CAS No.:480-41-1 Formula: C15H12O5 Molecular Weight : 272.26 . Synonyms: 4H-1-Benzopyran-4-one,2,3-dihydro-5,7-dihydroxy-2-(4-hydroxyphenyl)-,(57263395,S)-;Flavanone,4',5,7-trihydroxy-(8CI);Naringenin

HYDROXYBENZOIC ACID CAS 99-96-7

Name: 4-Hydroxybenzoic acid CAS No.:99-96-7 Formula: C7H6O3 Molecular Weight : 138.12 . Synonyms: p-Oxybenzoesaure [German];Benzoic acid, p-hydroxy-;Benzoic acid, p-hydroxy;4-Carboxyphenol;Benzoic acid, 4-hydroxy-, monosodium salt;4-Hydroxybenzoate;Hydroxybenzenecarboxylic acid;4-Hydroxybenzoesaeure;Benzoic acid,

3,4-Dimethoxy-2-pyridinemethanol CAS 72830-08-1

Name: 3,4-Dimethoxy-2-pyridinemethanol CAS No.:72830-08-1 Formula: C8H11NO3 Molecular Weight : 169.18 Synonyms: 2-Hydroxymethyl-3,4-dimethoxypyridine;3,4-Dimethoxy-2-(hydroxymethyl)pyridine; Density: 1.171 g/cm3 Boiling Point: 275.516 °C at 760 mmHg Flash Point: 120.428 °C

Homoharringtonine CAS 26833-87-4

Name: Homoharringtonine CAS No.:26833-87-4 Formula: C29H39NO9 Molecular Weight : 545.69 Synonyms: Cephalotaxine,4-methyl 2-hydroxy-2-(4-hydroxy-4-methylpentyl)butanedioate (ester), [3(R)]-;Homoharringtonine (8CI);CGX 635;Ceflatonin;Myelostat;NSC 141633;Omacetaxine mepesuccinate;Ephalotaxine,

PHYSOSTIGMINE SULFATE CAS 64-47-1

Name: PHYSOSTIGMINE SULFATE CAS No.:64-47-1 Formula: C30H42N6O4S Molecular Weight : 646.84 Synonyms: (3as-cis)-amate(estersulfate(2:1)(salt);(3as-cis)-ylcarbamate(estersulfate(2:1);3-b)indole,1,2,3,3a,8,8a-hexahydro-5-hydroxy-1,3a,8-trimethyl-pyrrolo(meth;eserinesulphate;physostigmineso4;physostigminesulphate;ESERINE

Glucosamine sulfate CAS 14999-43-0

Name: Glucosamine sulfate CAS No.:14999-43-0 Formula: 2(C6H13NO5).H2SO4 Molecular Weight : 456.42 Synonyms: Bis(2-ammonio-2-deoxy-D-glucose) sulphate;(2R,3R,4R,5S,6R)-3-Amino-6-(hydroxymethyl)oxane-2,4,5-triol; sulfuric acid;D-Glucose, 2-amino-2-deoxy-, sulfate (2:1); EINECS: 239-088-2 Boiling Point: 961.9 °C at 760

2-Propenoic acid,2-methyl-, 1,1-dimethylethyl ester CAS 585-07-9

Name: 2-Propenoic acid,2-methyl-, 1,1-dimethylethyl ester CAS No.:585-07-9 Formula: C8H14O2 Molecular Weight : 142.1956 Synonyms: Methacrylicacid, tert-butyl ester (6CI,7CI,8CI);1,1-Dimethylethyl methacrylate;AcryesterTB;Acryester TBMA;Light Ester TB;NSC 20957;tert-Butyl 2-methylacrylate;tert-Butyl

Betamethasone CAS 378-44-9

Name: Betamethasone CAS No.:378-44-9 Formula: C22H29FO5 Molecular Weight : 434.50 Synonyms: Celestene;Prestwick_703;9-fluoro-11,17-dihydroxy-17-(2-hydroxyacetyl)-10,13,16-trimethyl-6,7,8,11,12,14,15,16-octahydrocyclopenta[a]phenanthren-3-one;Rinderon A;Cidoten;beta-Methasone

3,4-Pyridinedicarboxylicacid, 5-hydroxy-, 4-methyl ester CAS 243980-03-2

Name: 3,4-Pyridinedicarboxylicacid, 5-hydroxy-, 4-methyl ester CAS No.:243980-03-2 Formula: C8H7NO5 Molecular Weight : 197.1449 Synonyms: 5-Hydroxypyridine-3,4-dicarboxylic acid 4-methyl ester;QC-9270;OR4210; Density: 1.499 g/cm3 Boiling Point: 476.2 °C at 760 mmHg Flash Point: 241.8 °C

1H-Imidazole,2,2'-[1,2-diazenediylbis(1-methylethylidene)]bis[4,5-dihydro-, hydrochloride(1:2) CAS 27776-21-2

Name: 1H-Imidazole,2,2'-[1,2-diazenediylbis(1-methylethylidene)]bis[4,5-dihydro-, hydrochloride(1:2) CAS No.:27776-21-2 Formula: C12H24Cl2N6 Molecular Weight : 322.14 Synonyms: AIBI;V 44;VA 044;Vazo 044;Vazo 44WSP;Wako VA 044; EINECS: 248-655-3 Boiling Point: 413.3 °C at 760 mmHg Flash Point: 203.7 °C

3-Pyridinecarbonylchloride, 5-bromo- CAS 39620-02-5

Name: 3-Pyridinecarbonylchloride, 5-bromo- CAS No.:39620-02-5 Formula: C6H3BrClNO Molecular Weight : 220.45 Synonyms: 5-Bromo-3-pyridinecarbonylchloride; Density: 1.76 g/cm3 Boiling Point: 265.4 °C at 760 mmHg Flash Point: 114.3 °C

D-Glucitol,1-deoxy-1-(methylamino)- CAS 6284-40-8

Name: D-Glucitol,1-deoxy-1-(methylamino)- CAS No.:6284-40-8 Formula: C7H17NO5 Molecular Weight : 195.22 Synonyms: Glucitol,1-deoxy-1-(methylamino)-, D- (8CI);Sorbitol, 1-deoxy-1-methylamino-

Thalidomide CAS 50-35-1

Name: Thalidomide CAS No.:50-35-1 Formula: C13H10N2O4 Molecular Weight : 258.25 Synonyms: Phthalimide,N-(2,6-dioxo-3-piperidyl)- (6CI,7CI,8CI);1,3-Dioxo-2-(2,6-dioxopiperidin-3-yl)isoindoline;3-Phthalimidoglutarimide;Celgene;Contergan;Distaval;K

Bismuth nitrate pentahydrate CAS 10035-06-0

Name: Bismuth nitrate pentahydrate CAS No.:10035-06-0 Formula: Bi(NO3)3.5(H2O) Molecular Weight : 485.07 Synonyms: Nitric acid,bismuth(3+) salt, pentahydrate (9CI);Bismuth nitrate (Bi(NO3)3) pentahydrate;Bismuth trinitrate pentahydrate;Bismuth(III) nitrate pentahydrate;Granions deBismuth; EINECS: 233-791-8 Density:

Piperazine,1-[(4-chlorophenyl)phenylmethyl]-4-[(3-methylphenyl)methyl]-, hydrochloride(1:2) CAS 1104-22-9

Name: Piperazine,1-[(4-chlorophenyl)phenylmethyl]-4-[(3-methylphenyl)methyl]-, hydrochloride(1:2) CAS No.:1104-22-9 Formula: C25H29Cl3N2 Molecular Weight : 463.91 Synonyms: Piperazine,1-(p-chloro-a-phenylbenzyl)-4-(m-methylbenzyl)-,dihydrochloride

Benzenebutanoic acid, a-amino-, ethyl ester,hydrochloride (1:1),(57263373,aR)- CAS 90940-54-8

Name: Benzenebutanoic acid, a-amino-, ethyl ester,hydrochloride (1:1),(57263374,aR)- CAS No.:90940-54-8 Formula: C12H17 N O2 . Cl H Molecular Weight : 243.72982 Synonyms: Benzenebutanoicacid, a-amino-, ethyl ester,hydrochloride,(57263375,R)-; Benzenebutanoic acid, a-amino-, ethyl ester, hydrochloride,(57263376,aR)-

3-Methoxypropylamine CAS 5332-73-0

Name: 3-Methoxypropylamine CAS No.:5332-73-0 Formula: C4H11NO Molecular Weight : 89.16 Synonyms: Propylamine,3-methoxy- (6CI,7CI,8CI);(3-Methoxypropan-1-yl)amine;1-Amino-3-methoxypropane;1-Methoxy-3-aminopropane;3-Amino-1-methoxypropane;3-Aminopropyl methyl

Androst-4-ene-3,17-dione,6-methylene- CAS 19457-55-7

Name: Androst-4-ene-3,17-dione,6-methylene- CAS No.:19457-55-7 Formula: C20H26O2 Molecular Weight : 298.42 Synonyms: 6-Methyleneandrost-4-en-3,17-dione; Density: 1.11 g/cm3 Boiling Point: 450.8 °C at 760 mmHg Flash Point: 168 °C

GCLE CAS 104146-10-3

Name: GCLE CAS No.:104146-10-3 Formula: C24H23ClN2O5S Molecular Weight : 486.9678 Synonyms: 5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid, 3-(chloromethyl)-8-oxo-7-[(phenylacetyl)amino]-,(57263366,4-methoxyphenyl)methylester,(57263367,6R,7R)- (9CI);5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylic

Azasetron Hydrochloride CAS 123040-69-7

Name: Azasetron Hydrochloride CAS No.:123040-69-7 Formula: C17H20ClN3O3 Molecular Weight : 386.28 Synonyms: UNII-77HC7URR9Z;(-

Ethyl phenylacetate CAS 101-97-3

Name: Ethyl phenylacetate CAS No.:101-97-3 Formula: C10H12O2 Molecular Weight : 164.20 Synonyms: Ethyl phenacetate;Benzeneacetic acid, ethyl ester;FEMA No. 2452;Phenylacetic acid, ethyl ester;Phenyl-acetic acid, ethyl ester;Acetic acid, phenyl-, ethyl ester;Ethyl .alpha.-toluate;alpha-Toluic acid, ethyl ester;Ethyl

2H-Oxacyclotetradecino[4,3-d]oxazole-2,6,8,14(1H,7H,9H)-tetrone,4-ethyloctahydro-11-methoxy-3a,7,9,11,13,15-hexamethyl-1-[4-[4-(3-pyridinyl)-1H-imidazol-1-yl]butyl]-10-[[3,4,6-trideoxy-3-(dimethylamino)-b-D-xylo-hexopyranosyl]oxy]-,(57263361,3aS,4R,7R,9R

Name: 2H-Oxacyclotetradecino[4,3-d]oxazole-2,6,8,14(1H,7H,9H)-tetrone,4-ethyloctahydro-11-methoxy-3a,7,9,11,13,15-hexamethyl-1-[4-[4-(3-pyridinyl)-1H-imidazol-1-yl]butyl]-10-[[3,4,6-trideoxy-3-(dimethylamino)-b-D-xylo-hexopyranosyl]oxy]-,(57263362,3aS,4R,7R,9R,10R,11R,13R,15R,15aR)- CAS No.:191114-48-4 Formula:

7-Xylosyltaxol CAS 90332-66-4

Name: 7-Xylosyltaxol CAS No.:90332-66-4 Formula: C52H59NO18 Molecular Weight : 986.03636 Synonyms: Density: 1.451 g/cm3 Boiling Point: 1062.337 °C at 760 mmHg Flash Point: 596.28 °C

Sitagliptin phosphate CAS 654671-78-0

Name: Sitagliptin phosphate CAS No.:654671-78-0 Formula: C16H18F6N5O5P Molecular Weight : 505.31 Synonyms: (2R)-4-Oxo-4-(3-(trifluoromethyl)-5,6-dihydro(1,2,4)triazolo(4,3-a)pyrazin-7(8H)-yl)-1-(2,4,5-trifluorophenyl)butan-2-amine phosphate salt;MK 0431; Boiling Point: 529.9 °C at 760 mmHg Flash Point: 274.3 °C

5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid, 7-[[(2Z)-2-(2-amino-4-thiazolyl)-2-(methoxyimino)acetyl]amino]-8-oxo-3-[[(1,2,5,6-tetrahydro-2-methyl-5,6-dioxo-1,2,4-triazin-3-yl)thio]methyl]-,sodium salt (1:2),(57263355,6R,7R)- CAS 74578-69-1

Name: 5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid, 7-[[(2Z)-2-(2-amino-4-thiazolyl)-2-(methoxyimino)acetyl]amino]-8-oxo-3-[[(1,2,5,6-tetrahydro-2-methyl-5,6-dioxo-1,2,4-triazin-3-yl)thio]methyl]-,sodium salt (1:2),(57263356,6R,7R)- CAS No.:74578-69-1 Formula: C18H18N8Na2O7S3 Molecular Weight :

Pipracil, Piper CAS 66258-76-2

Name: Pipracil, Piper CAS No.:66258-76-2 Formula: C23H29N5O8S Molecular Weight : 517.55 Synonyms:

Iodomethyl pivalate CAS 53064-79-2

Name: Iodomethyl pivalate CAS No.:53064-79-2 Formula: C6H11IO2 Molecular Weight : 242.06 Synonyms: IodomethylPivalate;LDO17:Iodomethyl Pivalate;iodomethyl 2,2-dimethylpropanoate; EINECS: 258-339-7 Density: 1.614 g/cm3 Boiling Point: 197.2 °C at 760 mmHg Flash Point: 73.1 °C Appearance: Sight yellow or colorless

Kinase(enzyme-activating), uro- CAS 9039-53-6

Name: Kinase(enzyme-activating), uro- CAS No.:9039-53-6 Formula: Unspecified Molecular Weight : 0 Synonyms: Abbokinase;Actosolv;Breokinase;Double-chainurokinase-type plasminogen activator;E.C. 3.4.21.31;E.C. 3.4.21.73;E.C.3.4.99.26;Persolv;Plasminokinase, urinary;Pro-HGF convertase;Pro-hepatocytegrowth factor

Benzoic acid,4-chloro-3,5-dinitro-, 2-methylpropyl ester CAS 58263-53-9

Name: Benzoic acid,4-chloro-3,5-dinitro-, 2-methylpropyl ester CAS No.:58263-53-9 Formula: C11H11ClN2O6 Molecular Weight : 302.67 Synonyms: 4-Chloro-3,5-dinitrobenzoicacid isobutyl ester; Density: 1.421 g/cm3 Boiling Point: 406.3 °C at 760 mmHg Flash Point: 199.5 °C

Myclobutanil CAS 88671-89-0

Name: Myclobutanil CAS No.:88671-89-0 Formula: C15H17ClN4 Molecular Weight : 288.78 Synonyms: Systhane;Nu-Flow M;Nova (pesticide);Rally;Systhane 6 Flo;RH 3866;1H-1,2,4-Triazole-1-propanenitrile,R-butyl-R-(4-chlorophenyl)-;NOVA

Atosiban CAS 90779-69-4

Name: Atosiban CAS No.:90779-69-4 Formula: C43H67N12O38S2 Molecular Weight : 994.19 Synonyms: Oxytocin,1-(3-mercaptopropanoicacid)-2-(O-ethyl-D-tyrosine)-4-L-threonine-8-L-ornithine-;Antocile;Antocin;Antocin II;CAP 449;CAP 476;CAP 581;F 314;ORF 22164;RW22164;RWJ 22164;Tractocil;Tractocile;Atosiban Acetate; Density:

4-(Trifluoromethyl)benzene-1-sulfonyl chloride CAS 2991-42-6

Name: 4-(Trifluoromethyl)benzene-1-sulfonyl chloride CAS No.:2991-42-6 Formula: C7H4ClF3O2S Molecular Weight : 244.61 Synonyms: p-Toluenesulfonylchloride, a,a,a-trifluoro- (6CI,7CI,8CI);[p-(Trifluoromethyl)phenyl]sulfonyl chloride;p-Trifluoromethylbenzenesulfonylchloride;a,a,a-Trifluoro-p-toluenesulfonyl chloride;

Manganese disodium EDTA trihydrate CAS 15375-84-5

Name: Manganese disodium EDTA trihydrate CAS No.:15375-84-5 Formula: C10H12N2O8MnNa2 Molecular Weight : 389.12 Synonyms: Disodium manganese ethylenediaminetetraacetate;Dissolvine E-Mn 13;EDTAdisodium manganese salt;Manganese disodium EDTA;Disodium manganeseEDTA;NSC 7345;Sequestrene NA 2MN;Manganate(2-),

2-Butanone,2,2',2''-[O,O',O''-(methylsilylidyne)trioxime] CAS 22984-54-9

Name: 2-Butanone,2,2',2''-[O,O',O''-(methylsilylidyne)trioxime] CAS No.:22984-54-9 Formula: C13H27N3O3Si Molecular Weight : 301.46 Synonyms: 2-Butanone,O,O',O''-(methylsilylidyne)trioxime

Erythromycin lactobionate CAS 3847-29-8

Name: Erythromycin lactobionate CAS No.:3847-29-8 Formula: C49H89NO25 Molecular Weight : 1092.22 Synonyms: Erythromycin lactobionate [JAN];1334-29-8;Erythromycin lactobionate (1:1) (salt);Lactobionic acid, compd. with erythromycin (1:1);Erythromycin. compd. with lactobionic acid;Erythrocin (TN);D-Gluconic acid,

Benzamide,N-[[(3,5-dichloro-2,4-difluorophenyl)amino]carbonyl]-2,6-difluoro- CAS 83121-18-0

Name: Benzamide,N-[[(3,5-dichloro-2,4-difluorophenyl)amino]carbonyl]-2,6-difluoro- CAS No.:83121-18-0 Formula: C14H6Cl2F4N2O2 Molecular Weight : 381.11 Synonyms: AC291898;CME 134;CME 13406;Calicide;HOE 522;MK 139;Nomolt;Nomolt agro;OMS

1,3-Diethylbenzene CAS 141-93-5

Name: 1,3-Diethylbenzene CAS No.:141-93-5 Formula: C10H14 Molecular Weight : 134.22 Synonyms: Benzene,m-diethyl- (8CI);1,3-Diethylbenzene;NSC 62102;m-Diethylbenzene; EINECS: 205-511-4 Density: 0.866 g/cm3 Boiling Point: 181.675 °C at 760 mmHg Flash Point: 50.556 °C

Amifostine CAS 20537-88-6

Name: Amifostine CAS No.:20537-88-6 Formula: C5H15N2O3PS Molecular Weight : 214.25 Synonyms: WR 2721;2-(3-aminopropylamino)ethylsulfanylphosphonic acid;Gammaphos;Aminopropylaminoethyl thiophosphate;Ethanethiol,2-[(3-aminopropyl)amino]-,dihydrogen phosphate (ester);Ethiofos;S-(2-(3-Aminopropylamino)ethyl)

Propanoic acid,2-methyl-, 1,1'-[2,2-dimethyl-1-(1-methylethyl)-1,3-propanediyl] ester CAS 6846-50-0

Name: Propanoic acid,2-methyl-, 1,1'-[2,2-dimethyl-1-(1-methylethyl)-1,3-propanediyl] ester CAS No.:6846-50-0 Formula: C16H30 O4 Molecular Weight : 286.41 Synonyms: Isobutyricacid, 1-isopropyl-2,2-dimethyltrimethylene ester (7CI,8CI);Propanoic acid,2-methyl-, 2,2-dimethyl-1-(1-methylethyl)-1,3-propanediyl ester

2(5H)-Furanone,5-hydroxy-4-methyl- CAS 40834-42-2

Name: 2(5H)-Furanone,5-hydroxy-4-methyl- CAS No.:40834-42-2 Formula: C5H6O3 Molecular Weight : 114.10 Synonyms: 4-Hydroxy-3-methylbut-2-en-4-olide;4-Methyl-5-hydroxy-2(5H)-furanone;5-Hydroxy-4-methyl-5H-furan-2-one;5-hydroxy-4-methylfuran-2(5H)-one;4-Methyl-5-hydroxy-2(5H)-furanone; Density: 1.35 g/cm3 Boiling Point:

Schizandrin B CAS 61281-37-6

Name: Schizandrin B CAS No.:61281-37-6 Formula: C23H28O6 Molecular Weight : 400.46 Synonyms: g-Schizandrin (7CI);Schisandrin B;Benzo(3,4)cycloocta(1,2-f)(1,3)benzodioxole, 5,6,7,8-tetrahydro-1,2,3,13-tetramethoxy-6,7-dimethyl-, stereoisomer;Wuweizisu

Methyl 2-hydroxyethylcellulose CAS 9032-42-2

Name: Methyl 2-hydroxyethylcellulose CAS No.:9032-42-2 Formula: C2H6O2.xCH4O.xUnspecified Molecular Weight : 0 Synonyms: Metolose SEB 02T;Tylose MN;Methyl hydroxyethylcellulose;SNB 100T;Metolose SEW 4000;Walocel MT 10.000GO;Walocel MKX 30.000PF01;Metolose SEB 30T;Metolose SE;Modocoll E 100;Hydroxyethyl methyl

1H-Imidazole, 2-methyl-5-nitro- CAS 88054-22-2

Name: 1H-Imidazole, 2-methyl-5-nitro- CAS No.:88054-22-2 Formula: C4H5N3O2 Molecular Weight :

Benzene,1,1'-(dimethoxysilylene)bis- CAS 6843-66-9

Name: Benzene,1,1'-(dimethoxysilylene)bis- CAS No.:6843-66-9 Formula: C14H16O2Si Molecular Weight : 244.36 Synonyms: Silane,dimethoxydiphenyl- (6CI,7CI,8CI,9CI);AY 43-047;AZ 6183;D 6010;Dimethoxydiphenylsilane;KBM 202;KBM 202LS5300;KBM202S;KBM 202SS;LS 5300;NSC 93509;OF 1 (silane);TSL 8172;Z 6074; EINECS:

Chondroitin 6-sulfate sodium salt CAS 12678-07-8

Name: Chondroitin 6-sulfate sodium salt CAS No.:12678-07-8 Formula: (C14H19NO14SNa2)n Molecular Weight : 1390.11000 Synonyms: Chondroitin sulfate C sodium salt;CSC-Na;Shark Chondroitin; Appearance: Off white to creamy powder

1-Cyclopropyl-6,7-difluoro-1,4-dihydro-8-methoxy-4-oxo-3-quinolinecarboxylic acid ethyl ester CAS 112811-71-9

Name: 1-Cyclopropyl-6,7-difluoro-1,4-dihydro-8-methoxy-4-oxo-3-quinolinecarboxylic acid ethyl ester CAS No.:112811-71-9 Formula: C16H15F2NO4 Molecular Weight :

1H,4H-[1,3]Thiazeto[3,2-a]quinoline-3-carboxylicacid, 6-fluoro-1-methyl-4-oxo-7-(1-piperazinyl)- CAS 112984-60-8

Name: 1H,4H-[1,3]Thiazeto[3,2-a]quinoline-3-carboxylicacid, 6-fluoro-1-methyl-4-oxo-7-(1-piperazinyl)- CAS No.:112984-60-8 Formula: C16H16FN3O3S Molecular Weight : 349.38 Synonyms: AF 3013;NAD 394;NM 394;6-Fluoro-1-methyl-4-oxo-7-(1-piperazinyl)-4H-[1,3]thiazeto[3,2-a]quinoline-3-carboxylic acid; Density: 1.58

5-TERT-BUTYL-2-HYDROXYBENZALDEHYDE CAS 2725-53-3

Name: Benzaldehyde,5-(1,1-dimethylethyl)-2-hydroxy- CAS No.:2725-53-3 Formula: C11H14O2 Molecular Weight : 178.23 Synonyms: Salicylaldehyde,5-tert-butyl- (6CI,7CI);2-Hydroxy-5-tert-butylbenzaldehyde;5-(1,1-Dimethylethyl)salicylaldehyde;5-tert-Butyl-2-hydroxybenzaldehyde;5-tert-Butylsalicylaldehyde; Density:

2-Cyclopentenone CAS 930-30-3

Name: 2-Cyclopentenone CAS No.:930-30-3 Formula: C5H6O Molecular Weight : 82.10 Synonyms: 1-Cyclopenten-3-one;1-Cyclopenten-5-one;2-Cyclopentenone;Cyclopenten-3-one;NSC 73117;2-Cyclopenten-1-one; EINECS: 213-213-0 Density: 1.046 g/cm3 Boiling Point: 135.999 °C at 760 mmHg Flash Point: 40.045 °C Solubility: Almost

Mosapride citrate CAS 112885-42-4

Name: Mosapride citrate CAS No.:112885-42-4 Formula: C21H25ClFN3O3.C6H8O7 Molecular Weight : 614.02 Synonyms: Gasmotin;Benzamide, 4-amino-5-chloro-2-ethoxy-N-((4-((4-fluorophenyl)methyl)-2-morpholinyl)methyl)-, 2-hydroxy-1,2,3-propanetricarboxylate (1:1);Benzamide,4-amino-5-chloro-2-ethoxy-N-

Leuprorelin CAS 53714-56-0

Name: Leuprorelin CAS No.:53714-56-0 Formula: C59H84N16O12 Molecular Weight : 1209.41 Synonyms: Luteinizinghormone-releasing factor (pig),6-D-leucine-9-(N-ethyl-L-prolinamide)-10-deglycinamide-;(D-Leu6,des-Gly-NH210)-LH-RH ethylamide;D-Leu6-des-Gly10-LH-releasing hormone ethylamide;Des-Gly10-[D-Leu6]-LH-releasing

Amlodipine besylate CAS 111470-99-6

Name: Amlodipine besylate CAS No.:111470-99-6 Formula: C20H25ClN2O5.C6H6O3S Molecular Weight : 567.05 Synonyms: Amlodipine benzenesulfonate;Monopina;Amlodin;Norvas;3,5-Pyridinedicarboxylic acid,2-[(2-aminoethoxy)methyl]-4-(2-chlorophenyl)- 1,4-dihydro-6-methyl-,3-ethyl 5-methyl

-Azetidineacetic acid,3-[(1R)-1-[[(1,1-dimethylethyl)dimethylsilyl]oxy]ethyl]-a-methyl-4-oxo-,(57263320,aR,2S,3S)- CAS 90776-58-2

Name: 2-Azetidineacetic acid,3-[(1R)-1-[[(1,1-dimethylethyl)dimethylsilyl]oxy]ethyl]-a-methyl-4-oxo-,(57263321,aR,2S,3S)- CAS No.:90776-58-2 Formula: C14H27NO4Si Molecular Weight : 301.45 Synonyms: 2-Azetidineaceticacid, 3-[1-[[(1,1-dimethylethyl)dimethylsilyl]oxy]ethyl]-a-methyl-4-oxo-,

tetrabromoethane CAS 25167-20-8

Name: 1,1,1,2-Tetrabromoethane CAS No.:25167-20-8 Formula: C2H2 Br4 Molecular Weight : 345.65 Synonyms: Tetrabromoethane EINECS: 246-687-2 Density: 2.967 g/mL at 25 °C(lit.) Boiling Point: 227.9 °C at 760 mmHg Flash Point: 91.9 °C Appearance: white or light yellow inflammable liquid

Rizatriptan Benzoate CAS 145202-66-0

Name: Rizatriptan benzoate CAS No.:145202-66-0 Formula: C15H19N5 Molecular Weight : 391.47 Synonyms: 1H-Indole-3-ethanamine,N,N-dimethyl-5-(1H-1,2,4-triazol-1-ylmethyl)-, monobenzoate (9CI);MK 462;Maxalt;1H-Indole-3-ethanamine,N,N-dimethyl-5-(1H-1,2,4-triazol-1-ylmethyl)-, benzoate (1:1);Rizatriptan Benzoate

3,5-Pyridinedicarboxylicacid, 2-[(2-aminoethoxy)methyl]-4-(2-chlorophenyl)-1,4-dihydro-6-methyl-,3-ethyl 5-methyl ester CAS 88150-42-9

Name: 3,5-Pyridinedicarboxylicacid, 2-[(2-aminoethoxy)methyl]-4-(2-chlorophenyl)-1,4-dihydro-6-methyl-,3-ethyl 5-methyl ester CAS No.:88150-42-9 Formula: C20H25ClN2O5 Molecular Weight : 408.88 Synonyms: (R,S)-Amlodipine;Amlopres;Intervask;Racemic Amlodipine;Amlodipine base;Amlodipine(base);Amlodipine; Density: 1.227

Acid Black 1 CAS 1064-48-8

Name: Acid Black 1 CAS No.:1064-48-8 Formula: C22H14N6Na2O9S2 Molecular Weight : 616.49 Synonyms: Buffalo Black NBR;Calcocid Blue Black;Concorde Leather Blue TB;Covalene Blue Black 6BN;C Black 1;D and C Black No. 1;Diacid Blue Black 10B;Dinacid Black 10BX;Eniacid Black SH;Erio Black HA;Eriosin Blue Black B;Everlan

Urea,N-[(4-chlorophenyl)methyl]-N-cyclopentyl-N'-phenyl- CAS 66063-05-6

Name: Urea,N-[(4-chlorophenyl)methyl]-N-cyclopentyl-N'-phenyl- CAS No.:66063-05-6 Formula: C19H21ClN2O Molecular Weight : 328.84 Synonyms: Monceraen250SC;Monceren;NTN 19701;Pencycuron; EINECS: 266-096-3 Density: 1.22 g/cm3 Boiling Point: 528.5 °C at 760 mmHg Flash Point: 273.4 °C

Roxithromycin CAS 80214-83-1

Name: Roxithromycin CAS No.:80214-83-1 Formula: C41H76N2O15 Molecular Weight : 837.05 Synonyms: Roxithromycin，Stearyl Glycyrrhetinate;Roxithromycin EP;RU 965;Rulide (TN);Roxithromycinum [Latin];RU 28965;9-(O-((2-Methoxyethoxy)methyl)oxime)erythromycin;Roxithromycin (JP14/USAN);Roxithromycine

p-Tolualdehyde CAS 104-87-0

Name: p-Tolualdehyde CAS No.:104-87-0 Formula: C8H8O Molecular Weight : 120.15 Synonyms: para-Toluyl aldehyde;p-Methyl benzaldehyde;4-Methyl benzaldehyde;p-tolualdehyde;4-methylBenzaldehyde;4-Toluicaldehyde;p-Formyltoluene;FEMA No. 3068;4-Toluylaldehyde;4-Methylbenzaldehyde;Benzaldehyde, 4-methyl-;Methylbenzaldehyde

Eudragit RL 12.5 CAS 33434-24-1

Name: Eudragit RL 12.5 CAS No.:33434-24-1 Formula: C19H34ClNO6 Molecular Weight : 407.9294 Synonyms: Density: g/cm3 Boiling Point: °Cat760mmHg Flash Point: °C

Vancocine hydrochloride CAS 1404-93-9

Name: Vancocine hydrochloride CAS No.:1404-93-9 Formula: C66H75Cl2N9O24.HCl Molecular Weight : 1485.86 Synonyms: Lyphocin;Meek;Vancocinehydrochloride;Vancocyn;Vancor; EINECS: 215-772-6 Density: 1.657 g/cm3 Solubility: 50 mg/mL in water,clear, yellow Appearance: off white powder

Propanedioic acid,2-(3-oxo-2-pentylcyclopentyl)-, 1,3-dimethyl ester CAS 51806-23-6

Name: Propanedioic acid,2-(3-oxo-2-pentylcyclopentyl)-, 1,3-dimethyl ester CAS No.:51806-23-6 Formula: C15H24 O5 Molecular Weight : 284.34806 Synonyms: Propanedioicacid,(57263304,3-oxo-2-pentylcyclopentyl)-, dimethyl ester (9CI); Dimethyl 2-n-pentyl-3-oxocyclopentylmalonate Density: 1.075g/cm3 Boiling Point:

Platinate(2-),hexachloro-, hydrogen, hydrate (1:2:6),(57263300,OC-6-11)- CAS 18497-13-7

Name: Platinate(2-),hexachloro-, hydrogen, hydrate (1:2:6),(57263301,OC-6-11)- CAS No.:18497-13-7 Formula: Cl6Pt . 6 H2 O . 2 H Molecular Weight : 521.97 Synonyms: Platinate(2-),hexachloro-, dihydrogen, hexahydrate (8CI); Platinate(2-), hexachloro-,dihydrogen, hexahydrate,(57263302,OC-6-11)- (9CI); Chloroplatinic

Tenatoprazole CAS 113712-98-4

Name: Tenatoprazole CAS No.:113712-98-4 Formula: C16H18N4O3S Molecular Weight : 346.4041 Synonyms: 1H-Imidazo[4,5-b]pyridine,5-methoxy-2-[[(4-methoxy-3,5-dimethyl-2-pyridinyl)methyl]sulfinyl]-

Polyethylene oxide CAS 25322-68-3

Name: Poly(ethylene oxide) CAS No.:25322-68-3 Formula: C2nH4n+2On+1 Molecular Weight : 0 Synonyms: Glycols, polyethylene;Glycols,polyethylene (8CI);AlkoxE 130;Alkox E 30;Alkox E 30G;Alkox R 100;Alkox R 15;Alkox R 400;Alkox SW;AquaCalk TWB-P;Aquaffin;Badimol;Breox 20M;Breox PEG 300;CBP 20;Carbowax 14000;Carbowax

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Roxatidine acetate hydrochloride CAS 93793-83-0

Name: Roxatidine acetate hydrochloride CAS No.:93793-83-0 Formula: C19H28N2O4.ClH Molecular Weight : 384.95 Synonyms: Altat (TN);Roxatidine acetate hydrochloride (JAN/USAN);Acetamide,2-(acetyloxy)-N-[3-[3-(1-piperidinylmethyl)phenoxy]propyl]-,monohydrochloride;Altat;Roxatidine aceacetate

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Barium hydroxide monohydrate CAS 22326-55-2

Name: Barium hydroxide monohydrate CAS No.:22326-55-2 Formula: Ba(OH)2.H2O Molecular Weight : 189.36 Synonyms: Bariumhydroxide, monohydrate (8CI);Barium hydroxide(Ba(OH)2), monohydrate (9CI); EINECS: 241-234-5 Density: 3.743 g/mL at 25 °C(lit.) Appearance: white powder

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Iodixanol CAS 92339-11-2

Name: Iodixanol CAS No.:92339-11-2 Formula: C35H44I6N6O15 Molecular Weight : 1550.18 Synonyms: Iodixanolum;OptiPrep;Visipaque;5,5'-((2-Hydroxytrimethylene)bis(acetylimino))bis(N,N'-bis(2,3-dihydroxypropyl)-2,4,6-triiodoisophthalamide);1,3-Benzenedicarboxamide,

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Benzeneacetic acid, a-(hydroxymethyl)-,(57263287,3-endo)-8-methyl-8-azabicyclo[3.2.1]oct-3-yl ester,(57263288,aS)- CAS 101-31-5

Name: Benzeneacetic acid, a-(hydroxymethyl)-,(57263289,3-endo)-8-methyl-8-azabicyclo[3.2.1]oct-3-yl ester,(57263290,aS)- CAS No.:101-31-5 Formula: C17H23NO3 Molecular Weight : 289.37 Synonyms: 1aH,5aH-Tropan-3a-ol,(57263291,-)-tropate (ester) (8CI);Benzeneacetic acid,

Nateglinide CAS 105816-04-4

Name: Nateglinide CAS No.:105816-04-4 Formula: C19H27NO3 Molecular Weight : 317.43 Synonyms: Nateglinide [INN];Fastic;trans-N-((4-(1-Methylethyl)cyclohexyl)carbonyl)-D-phenylalanine;A 4166;Starlix;Nateglinide(105816-04-4);Nateglinide,105816-04-4;Starsis

1,4-Benzoquinone CAS 106-51-4

Name: 1,4-Benzoquinone CAS No.:106-51-4 Formula: C6H4O2 Molecular Weight : 108.1 Synonyms: p-Benzoquinone(8CI);1,4-Cyclohexadienedione;Chinone;NSC 36324;PBQ 2;Stearer PBQ;p-Quinone;Quinone; EINECS: 203-405-2 Density: 1.256 g/cm3 Boiling Point: 174 ºC at 760 mmHg Flash Point: 59.3 ºC Solubility: 10 g/L (25 ºC) in

Eflornithine hydrochloride CAS 68278-23-9

Name: Eflornithine hydrochloride CAS No.:68278-23-9 Formula: C6H13ClF2N2O2 Molecular Weight : 218.63 Synonyms: DL-Ornithine,2-(difluoromethyl)-, monohydrochloride;Ornithine, 2-(difluoromethyl)-,monohydrochloride (9CI);NSC 270295;NSC 337250;Vaniqa; EINECS: 269-532-0 Boiling Point: 347 °C at 760 mmHg Flash Point: 163.7

Tungsten oxide (WO3) CAS 1314-35-8

Name: Tungsten oxide (WO3) CAS No.:1314-35-8 Formula: WO3 Molecular Weight : 231.85 Synonyms: C.I. 77901;Ditungsten hexaoxide;Tritungsten nonaoxide;Tungsten oxide;Tungsten oxide(W2O6);Tungsten oxide (W3O9);Tungsten oxide (W4O12);Tungsten trioxide (WO3);Tungsten(VI) oxide;Tungstic anhydride;Tungsticoxide;Tungstic oxide

Parthenolide CAS 20554-84-1

Name: Oxireno[9,10]cyclodeca[1,2-b]furan-9(1aH)-one,2,3,6,7,7a,8,10a,10b-octahydro-1a,5-dimethyl-8-methylene-,(57263281,1aR,4E,7aS,10aS,10bR)- CAS No.:20554-84-1 Formula: C15H20O3 Molecular Weight : 248.32 Synonyms: Germacra-1(10),11(13)-dien-12-oicacid, 4,5a-epoxy-6b-hydroxy-, g-lactone

Benzene,1-(chloromethyl)-2-methyl- CAS 552-45-4

Name: Benzene,1-(chloromethyl)-2-methyl- CAS No.:552-45-4 Formula: C8H9Cl Molecular Weight : 140.61 Synonyms: o-Xylene, a-chloro- (7CI,8CI);1-(Chloromethyl)-2-methylbenzene;1-Methyl-2-chloromethylbenzene;2-(Chloromethyl)toluene;o-Chloromethyltoluene;o-Methylbenzyl chloride;o-Xylyl

Acetamide,N-(4-ethoxyphenyl)- CAS 62-44-2

Name: Acetamide,N-(4-ethoxyphenyl)- CAS No.:62-44-2 Formula: C10H13NO2 Molecular Weight :

Disperse Orange 37 CAS 12223-33-5

Name: Disperse Orange 37 CAS No.:12223-33-5 Molecular Weight : 392.245 Synonyms: C.I. Disperse Orange 37 EINECS: 236-325-1 Density: 1.35g/cm3 Boiling Point: 603.1°C at 760 mmHg Flash Point: 318.6°C

1H-Inden-1-one,2,3-dihydro-5,6-dimethoxy-2-(4-pyridinylmethylene)- CAS 4803-74-1

Name: 1H-Inden-1-one,2,3-dihydro-5,6-dimethoxy-2-(4-pyridinylmethylene)- CAS No.:4803-74-1 Formula: C17H15NO3 Molecular Weight : 281.31 Synonyms: 1-Indanone,5,6-dimethoxy-2-(4-pyridylmethylene)-(7CI,8CI); Density: 1.256 g/cm3 Boiling Point: 499.9 °C at 760 mmHg Flash Point: 256.1 °C

Disulfide,2-propen-1-yl propyl CAS 2179-59-1

Name: Disulfide,2-propen-1-yl propyl CAS No.:2179-59-1 Formula: C6H12S2 Molecular Weight : 148.30 Synonyms: Disulfide,2-propenyl propyl (9CI);Disulfide, allyl propyl (6CI,7CI,8CI);2-Propenylpropyl disulfide;4,5-Dithia-1-octene;Propyl allyldisulfide; EINECS: 218-550-7 Density: 0.984 g/cm3 Boiling Point: 178.2 °C at 760

5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid,7-[[(2R)-2-[[(4-ethyl-2,3-dioxo-1-piperazinyl)carbonyl]amino]-2-(4-hydroxyphenyl)acetyl]amino]-3-[[(1-methyl-1H-tetrazol-5-yl)thio]methyl]-8-oxo-,sodium salt (1:1),(57263272,6R,7R)- CAS 62893-20-3

Name: 5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylicacid,7-[[(2R)-2-[[(4-ethyl-2,3-dioxo-1-piperazinyl)carbonyl]amino]-2-(4-hydroxyphenyl)acetyl]amino]-3-[[(1-methyl-1H-tetrazol-5-yl)thio]methyl]-8-oxo-,sodium salt (1:1),(57263273,6R,7R)- CAS No.:62893-20-3 Formula: C25H27N9O8S2. Na Molecular Weight :

Pyridinium,1-[[(6R,7R)-7-[[(2Z)-(2-amino-4-thiazolyl)[(1-carboxy-1-methylethoxy)imino]acetyl]amino]-2-carboxy-8-oxo-5-thia-1-azabicyclo[4.2.0]oct-2-en-3-yl]methyl]-,inner salt, hydrate (1:5) CAS 78439-06-2

Name: Pyridinium,1-[[(6R,7R)-7-[[(2Z)-(2-amino-4-thiazolyl)[(1-carboxy-1-methylethoxy)imino]acetyl]amino]-2-carboxy-8-oxo-5-thia-1-azabicyclo[4.2.0]oct-2-en-3-yl]methyl]-,inner salt, hydrate (1:5) CAS No.:78439-06-2 Formula: C22H22N6O7S2. 5H2O Molecular Weight :

1,3,4-Thiadiazole-2(3H)-thione,5-amino- CAS 2349-67-9

Name: 1,3,4-Thiadiazole-2(3H)-thione,5-amino- CAS No.:2349-67-9 Formula: C2H3N3S2 Molecular Weight : 133.20 Synonyms: 1,3,4-Thiadiazole-2-thiol,5-amino-

2,4(1H,3H)-Pyrimidinedione,6-[[3-[4-(2-methoxyphenyl)-1-piperazinyl]propyl]amino]-1,3-dimethyl- CAS 34661-75-1

Name: 2,4(1H,3H)-Pyrimidinedione,6-[[3-[4-(2-methoxyphenyl)-1-piperazinyl]propyl]amino]-1,3-dimethyl- CAS No.:34661-75-1 Formula: C20H29N5O3 Molecular Weight : 387.54 Synonyms: B 66256;BKU;Ebrantil;Eupressyl;Mediatensyl;NSC 310405;Uraprene;Uropidil; EINECS: 252-130-4 Density: 1.26 g/cm3 Boiling Point: 549 °C at 760

3-Octyn-1-ol CAS 14916-80-4

Name: 3-Octyn-1-ol CAS No.:14916-80-4 Formula: C8H14 O Molecular Weight : 126.20 . CODE Synonyms:NSC 67251 EINECS: 238-986-1 Density: 0.876 Boiling Point: 86 - 88 C at 9 mm Hg Flash Point: > 110 C Solubility: Soluble View

C.I.Direct Yellow 96 CAS 61725-08-4

Name: C.I.Direct Yellow 96 CAS No.:61725-08-4 Molecular Weight : 0 Synonyms: DiphenylBrilliant Flavine 7GFF; Solophenyl Flavine 7GFE; Solophenyl Flavine 7GFE 500

2-Oxazolidinone,5-(4-morpholinylmethyl)-3-[[(5-nitro-2-furanyl)methylene]amino]- CAS 139-91-3

Name: 2-Oxazolidinone,5-(4-morpholinylmethyl)-3-[[(5-nitro-2-furanyl)methylene]amino]- CAS No.:139-91-3 Formula: C13H16N4O6 Molecular Weight : 324.33 Synonyms: 2-Oxazolidinone,5-(morpholinomethyl)-3-[(5-nitrofurfurylidene)amino]-

Potassium sulfate CAS 7778-80-5

Name: Potassium sulfate CAS No.:7778-80-5 Formula: K2SO4 Molecular Weight : 174.26 Synonyms: Sulfuricacid dipotassium salt (8CI,9CI);Arcanum duplicatum;Dipotassium sulfate;Sulfuric acid potassiumsalt (1:2);Potassium sulfate (K2(SO4));Potassium sulfate (K2SO4);Potassium sulphate;Sal polychrestum;Sulfuric acid,

Platinum,diamminebis(nitrito-kN)- CAS 14286-02-3

Name: Platinum,diamminebis(nitrito-kN)- CAS No.:14286-02-3 Formula: H6N4O4Pt Molecular Weight : 321.15 Synonyms: Platinum,diamminebis(nitrito-N)-;Platinum, diamminedinitro- (8CI);Platinum,dinitrodiammine

(7CI);Diamminedinitroplatinum;NSC 130242; EINECS: 238-203-3 Density: 1.015 g/mL at 25 °C Appearance: colorless

Manganese bromide CAS 13446-03-2

Name: Manganese bromide CAS No.:13446-03-2 Formula: Br2Mn Molecular Weight : 214.75 Synonyms: Manganesebromide;Manganese dibromide;Manganese(II) bromide;Manganous bromide; EINECS: 236-591-9 Density: 4.385 g/mL at 25 °C(lit.) Appearance: Reddish crystalline powder

Diacetin CAS 25395-31-7

Name: Diacetin CAS No.:25395-31-7 Formula: C7H12O5 Molecular Weight : 176.19 Synonyms: 1,2,3-Propanetriol,esters,diacetate;Estol 1582;[(2S)-2-acetyloxy-3-hydroxy-propyl] acetate;Estol 1583;(3-acetyloxy-2-hydroxy-propyl) acetate;[(2R)-2-acetyloxy-3-hydroxy-propyl] acetate;Glycerol diacetate; EINECS: 246-941-2 Density:

4,4'-DIAMINODIPHENYLAMINE SULF CAS 537-65-5

Name: 1,4-Benzenediamine,N1-(4-aminophenyl)- CAS No.:537-65-5 Formula: C12H13 N3 Molecular Weight : 199.25 Synonyms: 1,4-Benzenediamine,N-(4-aminophenyl)- (9CI); Diphenylamine, 4,4'-diamino-(6CI,7CI,8CI);4,4'-Diaminodiphenylamine; 4,4'-Iminodianiline; Benzenamine, 4,4'-iminobis-;Bis(4-aminophenyl)amine;

1,8-Diazafluoran-9-one CAS 54078-29-4

Name: 1,8-Diazafluoran-9-one CAS No.:54078-29-4 Formula: C11H6N2O Molecular Weight : 182.18 Synonyms: 1,8-Diaza-9-fluorenone;1,8-Diazafluorenone;9H-1,8-Diazafluoren-9-one; Density: 1.387 g/cm3 Boiling Point: 396.3 °C at 760 mmHg Flash Point: 196.1 °C Appearance: yellow powder

Nelfinavir mesylate CAS 159989-65-8

Name: Nelfinavir mesylate CAS No.:159989-65-8 Formula: C32H45N3O4S.CH4O3S Molecular Weight : 663.89 Synonyms: Nelfinavir nesylate;[(3S-(3R*,4aR*,8aR*,2S*,3S*)]-2-[2-Hydroxy-3-phenylthiomethyl-4-aza-5-oxo-5-(2-methyl-3-hydroxy-phenyl)pentyl]-decahydroisoquinoline-3-N-t-butylcarboxamide methanesulfonic acid); Boiling

L-Lysine,N6-(1-oxododecyl)- CAS 52315-75-0

Name: L-Lysine,N6-(1-oxododecyl)- CAS No.:52315-75-0 Formula: C18H36N2O3 Molecular Weight : 328.49 Synonyms: Lysine,N6-lauroyl-, L- (6CI);Amihope LL;Famex L 12;N-Lauroyl-L-lysine;N6-Lauroyl-L-lysine;Ne-Lauroyl-L-lysine;Ne-Lauroyllysine;Ne-Lauroyllysine; EINECS: 257-843-4 Density: 0.994 g/cm3 Boiling Point: 528 °C at

Carbamodithioic acid,N,N-dimethyl-,sodium salt (1:1) CAS 128-04-1

Name: Carbamodithioic acid,N,N-dimethyl-,sodium salt (1:1) CAS No.:128-04-1 Formula: C3H7NNaS2 Molecular Weight : 144.21 Synonyms: (Dimethyldithiocarbamato)sodium;Dimethyldithiocarbamicacid sodium salt;N,N-Dimethyldithiocarbamic acid sodium salt; EINECS: 204-876-7 Density: 1.17 g/cm3 Boiling Point: 129.4 °C at 760

7-Oxabicyclo[4.1.0]heptane-3,4-dicarboxylicacid,3,4-bis(2-oxiranylmethyl) ester CAS 25293-64-5

Name: 7-Oxabicyclo[4.1.0]heptane-3,4-dicarboxylicacid,3,4-bis(2-oxiranylmethyl) ester CAS No.:25293-64-5 Formula: C14H18O7 Molecular Weight : 298.29 Synonyms: Bis(oxiran-2-ylmethyl) 7-oxabicyclo[4.1.0]heptane-3,4-dicarboxylate;1-Propanol, 2,3-epoxy-,7-oxabicyclo[4.1.0]heptane-3,4-dicarboxylate(2:1)(8CI);Diglycidyl

Vinpocetine CAS 42971-09-5

Name: Vinpocetine CAS No.:42971-09-5 Formula: C22H26N2O2 Molecular Weight : 350.50 Synonyms: (+)-Apovincaminic acid ethyl ester;(+)-Vinpocetine;(+)-cis-Apovincaminic acid ethyl ester;AY 27255;Apovincaminic acid ethyl ester;Bravinton;Cavinton;Ceractin;Ethyl(+)-apovincaminate;Ethyl (+)-cis-apovincaminate;Ethyl

Conjugated Estrogens CAS 12126-59-9

Name: ConjugatedEstrogens CAS No.:12126-59-9 Molecular Weight : 370.39500 Synonyms: Estrogens,conjugates;Ayerogen;Ayerogen Crema Vaginal;Azumon;C.E.S.;Cenestin;Climarest;Conjugated estrogens;Conjugates, estrogens;Conjugen;Dagynil;Emopremarin;Equin;Femavit;Hyphorin;Mannest;Menopak

Acrylic acid - maleic anhydride copolymer CAS 26677-99-6

Name: Acrylic acid - maleic anhydride copolymer CAS No.:26677-99-6 Formula: (C4H4O4)n.(C3H4O2)n Molecular Weight : 188.13500 Synonyms: MA/AA,MA/AA --Copolymer of Maleic and Acylic Acid;furan-2,5-dione; prop-2-enoic acid;Paama;2-Propenoic acid,polymers,polymer with 2,5-furandione;Copolymer of Maleic and Acylic Acid

Lamivudine CAS 134678-17-4

Name: Lamivudine CAS No.:134678-17-4 Formula: C8H11N3O3S Molecular Weight : 229.25 Synonyms: Epivir; GR109714X Density: 1.73 g/cm3 Boiling Point: 475.4 °C at 760mmHg Flash Point: 241.3 °C Appearance: white crystalline powder

1,2,3,9-Tetrahydro-4(H)-carbazol-4-one CAS 15128-52-6

Name: 1,2,3,9-Tetrahydro-4(H)-carbazol-4-one CAS No.:15128-52-6 Formula: C12H11NO Molecular Weight : 185.22 Synonyms: Carbazol-4(1H)-one,2,3-dihydro- (6CI,7CI,8CI);1,2,3,4-Tetrahydrocarbazol-4-one;1,2-Dihydro-4(3H)-carbazolone;2,3-Dihydrocarbazol-4(1H)-one;4-Oxo-1,2,3,4-tetrahydrocarbazole; Density: 1.275

Ethanone,1-(3-amino-2-hydroxyphenyl)- CAS 70977-72-9

Name: Ethanone,1-(3-amino-2-hydroxyphenyl)- CAS No.:70977-72-9 Formula: C8H9NO2 Molecular Weight : 151.16 Synonyms: Acetophenone,3'-amino-2'-hydroxy- (7CI);1-(3-Amino-2-hydroxyphenyl)ethanone;2-Acetyl-6-aminophenol;3-Amino-2-hydroxyacetophenone; Density: 1.242 g/cm3 Boiling Point: 287.2 °C at 760 mmHg Flash Point:

1-Azabicyclo[2.2.2]octan-3-one CAS 3731-38-2

Name: 1-Azabicyclo[2.2.2]octan-3-one CAS No.:3731-38-2 Formula: C7H11 N O Molecular Weight : 125.17 Synonyms: 3-Quinuclidinone(6CI,7CI,8CI);3-Oxoquinuclidine;3-Quinuclidone;NSC 168448; EINECS: 223-087-9 Density: 1.12g/cm3 Boiling Point: 204.9°Cat760mmHg Flash Point: 78.1°C

[1,1'-Biphenyl]-4-ol,4'-bromo- CAS 29558-77-8

Name: [1,1'-Biphenyl]-4-ol,4'-bromo- CAS No.:29558-77-8 Formula: C12H9BrO Molecular Weight : 249.11 Synonyms: 4-Biphenylol,4'-bromo- (8CI);Phenol, p-(p-bromophenyl)-

4-Bromobutyric acid CAS 2623-87-2

Name: 4-Bromobutyric acid CAS No.:2623-87-2 Formula: C4H7BrO2 Molecular Weight : 167.00 Synonyms:Butyric acid, 4-bromo-;4-bromobutanoate;Butanoic acid, 4-bromo-;NSC 99322; EINECS: 220-083-9 Density: 1.631 g/cm3 Boiling Point: 248.7 °C at 760 mmHg Flash Point: 111.3 °C Solubility: slightly soluble in water Appearance:

Benzenethiol,4,4'-thiobis- CAS 19362-77-7

Name: Benzenethiol,4,4'-thiobis- CAS No.:19362-77-7 Formula: C12H10S3 Molecular Weight : 250.39 Synonyms:Benzenethiol,4,4'-thiodi- (6CI,8CI);4,4'-Dimercaptodiphenyl sulfide;4,4'-Thiobis(phenylthiol);4,4'-Thiobisbenzenethiol;Bis(4-mercaptophenyl) sulfide;p,p'-Dimercaptodiphenyl sulfide; Density: 1.3 g/cm3 Boiling

1H-Tetrazole,5-(4-chlorobutyl)-1-cyclohexyl- CAS 73963-42-5

Name: 1H-Tetrazole,5-(4-chlorobutyl)-1-cyclohexyl- CAS No.:73963-42-5 Formula: C11H19ClN4 Molecular Weight : 242.75 Synonyms:1-Cyclohexyl-5-(4-chlorobutyl)-1,2,3,4-tetrazole;1-Cyclohexyl-5-(4-chlorobutyl)tetrazole; Density: 1.29 g/cm3 Boiling Point: 425.2 °C at 760 mmHg Flash Point: 210.9 °C Appearance: white

Acetamide, N-[5-(aminosulfonyl)-1,3,4-thiadiazol-2-yl]- CAS 59-66-5

Name: Acetamide, N-[5-(aminosulfonyl)-1,3,4-thiadiazol-2-yl]- CAS No.:59-66-5 Formula: C4H6N4O3S2 Molecular Weight : 222.26 Synonyms: Acetamide, N-(5-sulfamoyl-1,3,4-thiadiazol-2-yl)-

Gliclazide CAS 21187-98-4

Name: Gliclazide CAS No.:21187-98-4 Formula: C15H21N3O3S Molecular Weight : 323.41 Synonyms: Urea,1-(hexahydrocyclopenta[c]pyrrol-2(1H)-yl)-3-(p-tolylsulfonyl)- (8CI);Cyclopenta[c]pyrrole, benzenesulfonamide

Methyl dihydrojasmonate CAS 24851-98-7

Name: Methyl dihydrojasmonate CAS No.:24851-98-7 Formula: C13H22O3 Molecular Weight : 226.35 Synonyms: Cyclopentaneaceticacid, 3-oxo-2-pentyl-, methyl ester;Kharismal;Methyl (2-pentyl-3-oxocyclopentyl)acetate;Methyl(3-oxo-2-pentylcyclopentyl)acetate;methyl dihydrojasmonate (MDJ);Hedione; EINECS: 246-495-9 Density:

Pregnenolone CAS 145-13-1

Name: Pregnenolone CAS No.:145-13-1 Formula: C21H32O2 Molecular Weight : 316.48 Synonyms: 3-beta-Hydroxypregn-5-en-20-one;3beta-Hydroxy-Delta5-pregnen-20-one;3beta-Hydroxy-5-pregnene-20-one;Delta5-Pregnen-3beta-ol-20-one;Pregn-5-ene-3beta-ol-20-one;Pregnenolone:Pregna-5-ene-3beta-ol-20-one;Arthenolone; EINECS:

1H-Pyrazole,1-methyl- CAS 930-36-9

Name: 1H-Pyrazole,1-methyl- CAS No.:930-36-9 Formula: C4H6N2 Molecular Weight : 82.10 Synonyms:Pyrazole, 1-methyl- (6CI,7CI,8CI);1-Methyl-1H-pyrazole; EINECS: -0 Density: 0.996 g/cm3 Boiling

Point: 126.999 °C at 760 mmHg Flash Point: 26.222 °C Appearance: clear colorless to light yellow liquid Hazard Symbols: Xi

Zinc bacitracin CAS 1405-89-6

Name: Zinc bacitracin CAS No.:1405-89-6 Formula: C66H101N17O16SZn Molecular Weight : 1486.07 Synonyms: Baciferm(8CI);Albac;Bacifarmin;Bacifarmin-zinc;Baciferm PB;Noptracin;ST 56;Bacitracinzinc; EINECS: 215-787-8 Solubility: water: 5.1 g/L (28 °C) Appearance: White to pale buff powder

Ethyldiisopropylamine CAS 7087-68-5

Name: Ethyldiisopropylamine CAS No.:7087-68-5 Formula: C8H19N Molecular Weight : 129.24 Synonyms: 2-propanamine, N-ethyl-N-(1-methylethyl)-;N-ethyl-N-isopropylpropan-2-amine;2-Propanamine, N-ethyl-N- (1-methylethyl)-;1,1-Dimethyltriethylamine;N-Ethyl-N,

1,3-Bisbenzyl-2-oxoimidazolidine-4,5-dicarboxylic acid CAS 59564-78-2

Name: 1,3-Bisbenzyl-2-oxoimidazolidine-4,5-dicarboxylic acid CAS No.:59564-78-2 Formula: C19H18N2O5 Molecular Weight : 354.36 Synonyms: 1,3-Dibenzyl-2-oxoimidazolidine-4,5-dicarboxylic acid; EINECS: 261-806-8 Density: 1.426 g/cm3 Boiling Point: 654.2 °C at 760 mmHg Flash Point: 349.5 °C

Ketorolac CAS 74103-06-3

Name: Ketorolac CAS No.:74103-06-3 Formula: C15H13NO3 Molecular Weight : 255.27 Synonyms: 1H-Pyrrolizine-1-carboxylicacid, 5-benzoyl-2,3-dihydro-,(57263234,?à)-;1H-Pyrrolizine-1-carboxylicacid, 5-benzoyl-2,3-dihydro-;(+-)-5-Benzoyl-2,3-dihydro-1H-pyrrolizine-1-carboxylic acid;(+-)-Ketorolac; Density: 1.398

1,3-Propanediol,2-butyl-2-ethyl- CAS 115-84-4

Name: 1,3-Propanediol,2-butyl-2-ethyl- CAS No.:115-84-4 Formula: C9H20 O2 Molecular Weight : 160.26 Synonyms: 2-Butyl-2-ethyl-1,3-propanediol;2-Butyl-2-ethylpropanediol; 2-Ethyl-2-butyl-1,3-propanediol;3,3-Bis(hydroxymethyl)heptane; BEP; BEPD; DMH; NSC 406603; Nexcoat 700 EINECS: 204-111-7 Density: 0.93 g/cm3 Boiling

Benzenepropanoic acid, α-hydroxy-,(57263228,αR)- CAS 7326-19-4

Name: Benzenepropanoic acid, α-hydroxy-,(57263229,αR)- CAS No.:7326-19-4 Formula: C9H10O3 Molecular Weight : 166.17 Synonyms: Benzenepropanoicacid, a-hydroxy-,(57263230,R)-;Lactic acid,3-phenyl-,(57263231,R)- (8CI);(+)-2-Hydroxy-3-phenylpropanoic acid;(+)-3-Phenyllactic acid;(+)-b-Phenyllactic

Vinylidene chloride CAS 75-35-4

Name: Vinylidene chloride CAS No.:75-35-4 Formula: C2H2Cl2 Molecular Weight : 96.94 Synonyms: 1,1-Dichloroethene;1,1-Dichloroethylene;AI3-28804;Chlorure de vinylidene;Ethylene, 1,1-dichloro-;NCI-C54262;Vinylidene dichloride;as-Dichloroethylene;asym-Dichloroethylene; EINECS: 200-864-0 Density: 1.223 g/cm3 Boiling

2,3-Dichloropyrazine CAS 4858-85-9

Name: 2,3-Dichloropyrazine CAS No.:4858-85-9 Formula: C4H2Cl2N2 Molecular Weight : 148.98 Synonyms: Pyrazine,2,3-dichloro-; EINECS: 225-460-1 Density: 1.493 g/cm3 Boiling Point: 190.6 °C at 760 mmHg Flash Point: 86.1 °C Appearance: clear colorless to light yellow liquid Hazard Symbols: Xn, X

2-Bromo-5-fluorobenzoic acid CAS 394-28-5

Name: 2-Bromo-5-fluorobenzoic acid CAS No.:394-28-5 Formula: C7H4BrFO2 Molecular Weight : 219.01 Synonyms: 2-bromo-5-fluoro-benzoate;5-fluoro-2-bromobenzoic acid;2-Bromo-5-fluoro-benzoic acid; EINECS: -0 Density: 1.789 g/cm3 Boiling Point: 291.1 °C at 760 mmHg Flash Point: 129.8 °C Appearance: white to light yellow

1-Propanol,2-amino-3-(phenylmethoxy)-,(57263221,2R)- CAS 58577-87-0

Name: 1-Propanol,2-amino-3-(phenylmethoxy)-,(57263222,2R)- CAS No.:58577-87-0 Formula: C10H15 N O2 Molecular Weight : 181.23 Synonyms: 1-Propanol,2-amino-3-(phenylmethoxy)-,(57263223,R)-; (2R)-2-Amino-3-(benzyloxy)propan-1-ol;(R)-(+)-2-Amino-3-benzyloxy-1-propanol; (R)-2-Amino-3-benzyloxy-1-propanol Density:

Zinc,bis[N,N-bis(2-methylpropyl)carbamodithioato-kS,kS']-,(57263218,T-4)- CAS 36190-62-2

Name: Zinc,bis[N,N-bis(2-methylpropyl)carbamodithioato-kS,kS']-,(57263219,T-4)- CAS No.:36190-62-2 Formula: Zn.(C9H18NS2)2 Molecular Weight : 474.15 Synonyms: Carbamodithioicacid, bis(2-methylpropyl)-, zinc salt;Carbamodithioic acid,bis(2-methylpropyl)-, zinc complex;Bis(diisobutyldithiocarbamato)zinc;Isobutyl

2-Bromo-5-fluorobenzoic acid CAS 394-28-5

Name: 2-Bromo-5-fluorobenzoic acid CAS No.:394-28-5 Formula: C7H4BrFO2 Molecular Weight : 219.01 Synonyms: 2-bromo-5-fluoro-benzoate;5-fluoro-2-bromobenzoic acid;2-Bromo-5-fluoro-benzoic acid; EINECS: -0 Density: 1.789 g/cm3 Boiling Point: 291.1 °C at 760 mmHg Flash Point: 129.8 °C Appearance: white to light yellow

Irbesartan CAS 138402-11-6

Name: Irbesartan CAS No.:138402-11-6 Formula: C25H28N6O Molecular Weight : 428.54 Synonyms: BMS 186295;SR 47436;2-Butyl-3-(p-(o-1H-tetrazol-5-ylphenyl)benzyl)-1,3-diazaspiro(4.4)non-1-en-4-one;1,3-Diazaspiro[4.4]non-1-en-4-one,2-butyl- 3-[[2'-(1H-tetrazol-5-yl)[1,1'-biphenyl]-4-yl]-

Oleanic acid CAS 508-02-1

Name: Oleanic acid CAS No.:508-02-1 Formula: C30H48O3 Molecular Weight : 456.71 Synonyms: 3-beta-Hydroxyolean-12-en-28-oic acid;Olean-12-en-28-oic acid, 3-beta-hydroxy-

1-Pentanol, 5-amino- CAS 2508-29-4

Name: 1-Pentanol, 5-amino- CAS No.:2508-29-4 Formula: C5H13NO Molecular Weight : 103.16 Synonyms: 1-Amino-5-hydroxypentane;5-Aminopentanol;5-Hydroxy-1-pentanamine;5-Hydroxypentylamine;Pentanolamine; EINECS: 219-718-2 Density: 0.919 g/cm3 Boiling Point: 222.4 °C at 760 mmHg Flash Point: 65.6 °C Solubility: miscible in

Gemcitabine CAS 95058-81-4

Name: Gemcitabine CAS No.:95058-81-4 Formula: C9H11F2N3O4 Molecular Weight : 263.23 Synonyms: 4-amino-1-[(2R,4R,5R)-3,3-difluoro-4-hydroxy-5-(hydroxymethyl)oxolan-2-yl]pyrimidin-2-one;Cytidine, 2-deoxy-2,2-difluoro-2-Deoxy-.beta.-D-2,2-difluorocytidine;Gemcitabine [USAN:BAN:INN];DDFC;LY 188011;LY188011

2,2-Dimethoxyethylamine CAS 22483-09-6

Name: 2,2-Dimethoxyethylamine CAS No.:22483-09-6 Formula: C4H11NO2 Molecular Weight : 105.14 Synonyms: Acetaldehyde,amino-, dimethyl acetal (6CI,8CI);Ethylamine, 2,2-dimethoxy- (8CI);1-Amino-2,2-dimethoxyethane;Aminoacetaldehyde dimethyl acetal; EINECS: 245-026-5 Density: 0.944 g/cm3 Boiling Point: 133 °C at 760

Daidzin CAS 552-66-9

Name: Daidzin CAS No.:552-66-9 Formula: C21H20O9 Molecular Weight : 416.38 Synonyms: 7-(beta-D-Glucopyranosyloxy)-3-(4-hydroxyphenyl)-4H-1-benzopyran-4-one;Daidzein 7-glucoside;7-(beta-D-glucopyranosyloxy)-3-(4-hydroxyphenyl)-4H-chromen-4-one;4H-1-Benzopyran-4-one,7-(a-Dglucopyranosyloxy)-

Direct Green 26 CAS 6388-26-7

Name: Direct Green 26 CAS No.:6388-26-7 Formula: C50H33N12Na5O18S4 Molecular Weight :

Triethylamine CAS 121-44-8

Name: Triethylamine CAS No.:121-44-8 Formula: C6H15N Molecular Weight : 101.19 Synonyms: Triethylamine(7CI,8CI);(Diethylamino)ethane;N,N-Diethylethanamine;AI3-15425;CCRIS 4881;Ethanamine, N,N-diethyl-;HSDB 896;Triaethylamin;Trietilamina; EINECS: 204-469-4 Density: 0.75 g/cm3 Boiling Point: 90.495 °C at 760 mmHg Flash

Calcifediol anhydrous CAS 19356-17-3

Name: Calcifediol anhydrous CAS No.:19356-17-3 Formula: C27H44O2 Molecular Weight : 400.64 Synonyms: 9,10-Secocholesta-5,7,10(19)-triene-3,25-diol,(57263204,3b,5Z,7E)- (9CI);9,10-Secocholesta-5,7,10(19)-triene-3b,25-diol (8CI);25-HCC;25-Hydroxycholecalciferol;25-Hydroxyvitamin D;25-Hydroxyvitamin

3-(3,5-Dimethylphenoxy)propane-1,2-diol CAS 59365-66-1

Name: 3-(3,5-Dimethylphenoxy)propane-1,2-diol CAS No.:59365-66-1 Formula: C11H16O3 Molecular Weight : 196.24 Synonyms: 1,2-propanediol, 3-(3,5-dimethylphenoxy)-; Density: 1.125 g/cm3 Boiling Point: 372.7 °C at 760 mmHg Flash Point: 179.2 °C

Chlorthalidone CAS 77-36-1

Name: Benzenesulfonamide,2-chloro-5-(2,3-dihydro-1-hydroxy-3-oxo-1H-isoindol-1-yl)- CAS No.:77-36-1 Formula: C14H11ClN2O4S Molecular Weight : 338.78 Synonyms: Benzenesulfonamide,2-chloro-5-(1-hydroxy-3-oxo-1-isoindolinyl)-

2-Amino-3-nitropyridine CAS 4214-75-9

Name: 2-Amino-3-nitropyridine CAS No.:4214-75-9 Formula: C5H5N3O2 Molecular Weight : 139.11 Synonyms: 2-Pyridinamine, 3-nitro-;Pyridine, 2-amino-3-nitro-;5-nitro-1H-pyridin-6-amine;3-nitropyridin-2-amine;3-Nitro-2-aminopyridine;3-Nitropyridin-2-ylamine; EINECS: 224-144-0 Density: 1.437 g/cm3 Boiling Point: 299.1 °C at

Benzenamine,N-hydroxy-N-nitroso-, ammonium salt (1:1) CAS 135-20-6

Name: Benzenamine,N-hydroxy-N-nitroso-, ammonium salt (1:1) CAS No.:135-20-6 Formula: C6H9N3O2 Molecular Weight : 156.19 Synonyms: Benzenamine,N-hydroxy-N-nitroso-, ammonium salt (9CI);Hydroxylamine, N-nitroso-N-phenyl-,ammonium salt (8CI);Ammonium

3-Isothiazolecarboxylicacid, 5-nitro- CAS 36778-15-1

Name: 3-Isothiazolecarboxylicacid, 5-nitro- CAS No.:36778-15-1 Formula: C4H2 N2 O4 S Molecular Weight : 0 Synonyms: 5-Nitroisothiazole-3-carboxylicacid Density: 1.804g/cm3 Boiling Point: 253.795°C at 760 mmHg Flash Point: 107.292°C

Phenol, 2-(2-hydroxyethoxy)- CAS 4792-78-3

Name: Phenol, 2-(2-hydroxyethoxy)- CAS No.:4792-78-3 Formula: C8H10O3 Molecular Weight : 154.16 Synonyms: Ethanol,2-(o-hydroxyphenoxy)- (7CI,8CI);2-(2-Hydroxyethoxy)phenol;2-(2-Hydroxyphenoxy)ethanol;2-(2'-Hydroxyethoxy)phenol;2-(o-Hydroxyphenoxy)ethanol;2-[(2-Hydroxyphenyl)oxy]ethanol; EINECS: 225-346-1 Density:

R-(+)-alpha-Lipoic acid CAS 1200-22-2

Name: R-(+)-alpha-Lipoic acid CAS No.:1200-22-2 Formula: C8H14O2S2 Molecular Weight : 206.3256 Synonyms: 1,2-Dithiolane-3-pentanoicacid,(57263194,R)-;1,2-Dithiolane-3-valeric acid,(57263195,+)-(8CI);1,2-Dithiolane-3-pentanoicacid,(57263196,3R)-;(R)-Lipoic acid;(R)-a-Lipoic acid;Berlition;Byodinoral

1H-Indole-3-aceticacid, 5-methoxy- CAS 3471-31-6

Name: 1H-Indole-3-aceticacid, 5-methoxy- CAS No.:3471-31-6 Formula: C11H11NO3 Molecular Weight : 205.23 Synonyms: Indole-3-aceticacid, 5-methoxy- (6CI,7CI,8CI);(5-Methoxy-1H-indol-3-yl)acetic acid;2-(5-Methoxy-1H-indol-3-yl)acetic acid;2-(5-Methoxy-3-indolyl)acetic acid;5-Methoxy-1H-indole-3-acetic

4-tert-Butylcatechol CAS 98-29-3

Name: 4-tert-Butylcatechol CAS No.:98-29-3 Formula: C10H14O2 Molecular Weight : 166.22 Synonyms: p-t-Butylpyrocatechol;4-tert-Butyl-1,2-dihydroxybenzene;4-06-00-06014 (Beilstein Handbook Reference);4-tert-Butylpyrokatechin [Czech];p-tert-Butyl catechol;Synox TBC;Pyrocatechol,

1,4-Phenylene diisocyanate CAS 104-49-4

Name: 1,4-Phenylene diisocyanate CAS No.:104-49-4 Formula: C8H4N2O2 Molecular Weight : 160.14 Synonyms: 1,4-Diisocyanatobenzene;benzene, 1,4-diisocyanato-; EINECS: 203-207-6 Density: 1.17 g/cm3 Boiling Point: 236.2 °C at 760 mmHg Flash Point: 92.4 °C Appearance: white to light yellow crystals Hazard Symbols: Xn

L-Valine methyl ester hydrochloride CAS 6306-52-1

Name: L-Valine methyl ester hydrochloride CAS No.:6306-52-1 Formula: C6H14ClNO2 Molecular Weight : 167.64 Synonyms: H-Val-OMe·HCl; EINECS: 228-620-9 Boiling Point: 145.7 °C at 760 mmHg Flash Point: 20.7 °C Appearance: Crystalline Hazard Symbols: Xi

Meglumine Antimonate CAS 133-51-7

Name: Methylglucamine antimonate CAS No.:133-51-7 Formula: C7H17NO5.HSbO3 Molecular Weight : 365.98 Synonyms: Glucitol, 1-deoxy-1-(methylamino)-, compd. with antimonic acid (1:1), D-;Glucantime;Protostib;1-Deoxy-1-(methylamino)-D-glucitol, compound with antimonic acid (1:1);hydroxy-dioxo-stiborane;

L-Valine,N-[(2'-cyano[1,1'-biphenyl]-4-yl)methyl]-, methyl ester CAS 137863-89-9

Name: L-Valine,N-[(2'-cyano[1,1'-biphenyl]-4-yl)methyl]-, methyl ester CAS No.:137863-89-9 Formula: C20H22N2O2 Molecular Weight : 322.41 Synonyms: Methyl N-[(2'-cyano-4-biphenylyl)methyl]-L-valinate; Density: 1.12 g/cm3 Boiling Point: 477.9 °C at 760 mmHg Flash Point: 242.8 °C

Benzoic acid,5-acetyl-2-hydroxy-, methyl ester CAS 16475-90-4

Name: Benzoic acid,5-acetyl-2-hydroxy-, methyl ester CAS No.:16475-90-4 Formula: C10H10O4 Molecular Weight : 194.19 Synonyms: 5-Acetyl-2-hydroxybenzoic acidmethyl ester;Salicylicacid, 5-acetyl-, methyl ester (6CI,7CI,8CI);Methyl 3-acetyl-6-hydroxybenzoate;Methyl5-acetyl-2-hydroxybenzoate;NSC

Phenol, 3-mercapto- CAS 40248-84-8

Name: Phenol, 3-mercapto- CAS No.:40248-84-8 Formula: C6H6OS Molecular Weight : 126.18 Synonyms: Monothioresorcinol;m-Hydroxybenzenethiol;m-Mercaptophenol;3-Mercaptophenol;3-Hydroxythiophenol;3-Hydroxybenzenethiol;3-Hydroxyphenylmercaptan; Density: 1.255 g/cm3 Boiling Point: 256.3 °C at 760 mmHg Flash Point: 108.8

Phosphonitrilic chloride trimer CAS 940-71-6

Name: Phosphonitrilic chloride trimer CAS No.:940-71-6 Formula: Cl6N3P3 Molecular Weight : 347.66 Synonyms: 1,3,5,2,4,6-Triazatriphosphorine,2,2,4,4,6,6-hexachloro-2,2,4,4,6,6-hexahydro-(8CI,9CI);Phosphonitrilechloride, trimer

Disodium 5,5'-azodisalicylate CAS 6054-98-4

Name: Disodium 5,5'-azodisalicylate CAS No.:6054-98-4 Formula: C14H8N2Na2O6 Molecular Weight : 346.2027 Synonyms: Benzoicacid, 3,3'-azobis[6-hydroxy-, disodium salt (9CI);C.I. Mordant Yellow 5 (7CI);C.I. Mordant Yellow 5, disodium salt (8CI);3,3'-Azobis[6-hydroxybenzoic acid]disodium salt;Acid Chrome Flavine

Phenol,2-methoxy-4-methyl- CAS 93-51-6

Name: Phenol,2-methoxy-4-methyl- CAS No.:93-51-6 Formula: C8H10O2 Molecular Weight : 138.16 . Synonyms: Creosol(6CI);p-Cresol, 2-methoxy- (8CI);2-Methoxy-p-cresol;3-Methoxy-4-hydroxytoluene;4-Hydroxy-3-methoxy-1-methylbenzene;4-Hydroxy-3-methoxytoluene;4-Methyl-2-methoxyphenol;4-Methylguaiacol;Homoguaiacol;NSC

1,4-Benzenediol,2,5-dibromo- CAS 14753-51-6

Name: 1,4-Benzenediol,2,5-dibromo- CAS No.:14753-51-6 Formula: C6H4 Br2 O2 Molecular Weight : 267.9 Synonyms: Hydroquinone,2,5-dibromo- (6CI,8CI); 1,4-Dibromo-2,5-dihydroxybenzene;2,5-Dibromo-1,4-benzenediol; 2,5-Dibromo-1,4-hydroquinone;2,5-Dibromohydroquinone; NSC 133363 Density: 2.257g/cm3 Boiling Point:

2,5-Dibromopyridine CAS 624-28-2

Name: 2,5-Dibromopyridine CAS No.:624-28-2 Formula: C5H3Br2N Molecular Weight : 236.89 Synonyms: Pyridine, 2,5-dibromo-;2,5-Dibromo Pyridine; EINECS: 210-839-6 Density: 2.059 g/cm3 Boiling Point: 235.7 °C at 760 mmHg Flash Point: 96.4 °C Solubility: insoluble in water Appearance: Off-white crystals Hazard Symbols: Xn,

4-Methylmorpholine CAS 109-02-4

Name: 4-Methylmorpholine CAS No.:109-02-4 Formula: C5H11NO Molecular Weight : 101.15 Synonyms: 1-Methylmorpholine;4-Methylmorfolin;4-Methylmorfolin [Czech];Morpholine, N-methyl-;N-Methylmorpholine; EINECS: 203-640-0 Density: 0.931 g/cm3 Boiling Point: 114.1 °C at 760 mmHg Flash Point: 23.9 °C Solubility: Soluble

lactulose CAS 4618-18-2

Name: D-Fructose, 4-O-b-D-galactopyranosyl- CAS No.:4618-18-2 Formula: C12H22O11 Molecular Weight : 342.30 Synonyms: MLS-50;Piarle;Enulose;Duphalac;Liforos;Lagnos;Lactulose;Cholac;cephulac;laevolac;Caloryl; EINECS: 225-027-7 Density: 1.77 g/cm3 Boiling Point: 680.5 °C at 760 mmHg Flash Point: 365.4 °C Solubility: 76.4

Ciprofloxacin Hcl CAS 86483-48-9

Name: 3-Quinolinecarboxylicacid, 1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxo-7-(1-piperazinyl)-,hydrochloride (1:?) CAS No.:86483-48-9 Formula: C17H19ClFN3O3 Molecular Weight : 367.80 Synonyms: 3-Quinolinecarboxylicacid, 1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxo-7-(1-piperazinyl)-,hydrochloride

Benzalkonium Chloride CAS 85409-22-9

ame: Lutensit K-LC CAS No.:85409-22-9 Molecular Weight : 283.88000 Synonyms: Protectol KLC 50;Quaternary ammonium compounds, benzyl-C12-14-alkyldimethyl, chlorides;B 50 (surgactant);BAS 092-00E;Protectol KLC;Benzyl-C12-14-alkyldimethylammonium chlorides;Alkylbenzyldimethylammonium chlorides,

Quetiapine CAS 111974-69-7

Name: Quetiapine CAS No.:111974-69-7 Formula: C21H25N3O2S Molecular Weight : 383.51 Synonyms: 2-(2-(4-Dibenzo(b,f)(1,4)thiazepin-11-yl-1-piperazinyl)ethoxy)ethanol;Co-Quetiapine;Seroquel;Ethanol,2-[2-(4-dibenzo[b,f][1,4]thiazepin-11-yl-1-piperazinyl)ethoxy]-; Density: 1.27 g/cm3 Boiling Point: 556.5 °C at 760

Abacavir CAS 136470-78-5

Name: Abacavir CAS No.:136470-78-5 Formula: C14H18N6O Molecular Weight : 286.34 Synonyms: {(1S,4R)-4-[2-amino-6-(cyclopropylamino)-9H-purin-9-yl]cyclopent-2-en-1-yl}methanol;Ziagen;2-Cyclopentene-1-methanol, 4-(2-amino-6-(cyclopropylamino)-9H-purin-9-yl)-,(57263137,1S-cis)-;4-Amino-5-chloro-2-ethoxy benzoic

Aluminum bismuth oxide (Al3BiO6) CAS 12284-76-3

Name: Aluminum bismuth oxide (Al3BiO6) CAS No.:12284-76-3 Formula: Al3BiO6 Molecular Weight : 388.94500 Synonyms: Diwismut-tris(tetraoxoaluminat);UNII-JB5Y63JDHJ;Wismut-aluminat;Trialuminium bismuth hexaoxide; EINECS: 235-552-3

1,4-Dichlorobenzene CAS 106-46-7

Name: 1,4-Dichlorobenzene CAS No.:106-46-7 Formula: C6H4Cl2 Molecular Weight : 147.00 Synonyms: 1, 4-Dichlorobenzene;Paramoth;p-Dichloorbenzeen [Dutch];p-Diclorobenzene [Italian];p-Dichloorbenzeen;Dichlorobenzene, p-, solid;Benzene, p-dichloro-;1,4-Diclorobenzene [Italian];Paradichlorobenzol;Santochlor;RCRA waste no.

Pyridine,2-chloro-6-(trichloromethyl)- CAS 1929-82-4

Name: Pyridine,2-chloro-6-(trichloromethyl)- CAS No.:1929-82-4 Formula: C6H3Cl4N Molecular Weight : 230.91 Synonyms: 2-Chloro-6-Trichloromethylpyridine;2-Trichloromethyl-6-chloropyridine;6-Chloro-2-(trichloromethyl)pyridine;N-Serve;N-Serve 24E;2-Chloro-6-(trichloromethyl)pyridine; EINECS: 217-682-2 Density: 1.582

Vinpocetine CAS 42971-09-5

Name: Vinpocetine CAS No.:42971-09-5 Formula: C22H26N2O2 Molecular Weight : 350.50 Synonyms: (+)-Apovincaminic acid ethyl ester;(+)-Vinpocetine;(+)-cis-Apovincaminic acid ethyl ester;AY 27255;Apovincaminic acid ethyl ester;Bravinton;Cavinton;Ceractin;Ethyl(+)-apovincaminate;Ethyl (+)-cis-apovincaminate;Ethyl

4-Piperidinopiperidine CAS 4897-50-1

Name: 4-Piperidinopiperidine CAS No.:4897-50-1 Formula: C10H20N2 Molecular Weight : 168.2792 Synonyms: 1,4-Bipiperidyl;1,4'-bipiperidine;1-(3,4,5,6-tetrahydro-2H-pyridin-4-yl)-3,4,5,6-tetrahydro-2H-pyridine;4-(1-piperidinyl)piperidine;[1,4']Bipiperidinyl;4-(1-Piperidino)piperidine;1-(4-piperidyl)piperidine; EINECS:

1-(2-Nitrovinyl)benzene CAS 5153-67-3

Name: 1-(2-Nitrovinyl)benzene CAS No.:5153-67-3 Formula: C8H7NO2 Molecular Weight : 149.1467 Synonyms: Benzene,(57263127,2-nitroethenyl)-,(57263128,E)-;Styrene, b-nitro-,(57263129,E)-

DEET CAS 134-62-3

Name: N,N-Diethyl-3-methylbenzamide CAS No.:134-62-3 Formula: C12H17NO Molecular Weight : 191.30 Synonyms: m-Toluamide,N,N-diethyl- (6CI,7CI,8CI);3-Methyl-N,N-diethylbenzamide;AminceneC 140;Cutter Unscented;DETA;Delphene;Dieltamid;Diethyl-m-toluamide;Diethyltoluamide;ENT

Copper(I) cyanide CAS 544-92-3

Name: Copper(I) cyanide CAS No.:544-92-3 Formula: CuCN Molecular Weight : 89.56 Synonyms: Copper cyanide (Cu(CN));AI3-28745;Copper cyanide;Copper(+1) cyanide;Cupricin;Cuprous cyanide;HSDB 1438;RCRA waste number P029; EINECS: 208-883-6 Density: 2.92 g/mL at 25 °C(lit.) Boiling Point: 25.7 °C at 760 mmHg Solubility:

2,4(1H,3H)-Pyrimidinedione,1-b-D-arabinofuranosyl- CAS 3083-77-0

Name: 2,4(1H,3H)-Pyrimidinedione,1-b-D-arabinofuranosyl- CAS No.:3083-77-0 Formula: C9H12N2O6 Molecular Weight : 244.20 Synonyms: 1-[(2R,3S,4S,5R)-3,4-dihydroxy-5-(hydroxymethyl)oxolan-2-yl]pyrimidine-2,4-dione;1-pentofuranosylpyrimidine-2,4(1H,3H)-dione;Ara-U;Arabinofuranosyluracil;Arabinosyluracil;Arauridine;Uracil

Propanoic acid,2-formyl-3-oxo-, ethyl ester CAS 80370-42-9

Name: Propanoic acid,2-formyl-3-oxo-, ethyl ester CAS No.:80370-42-9 Formula: C6H8O4 Molecular Weight : 144.13 Synonyms: (Ethoxycarbonyl)malonaldehyde;2-Formyl-3-oxopropanoic acid ethylester;Ethyl 2,2-diformylacetate;Ethyl 2-formyl-3-oxopropanoate; Density: 1.143 g/cm3 Boiling Point: 181.475 °C at 760 mmHg Flash

4H-1-Benzopyran-4-one,3,7-dihydroxy-2-(3,4,5-trihydroxyphenyl)- CAS 490-31-3

Name: 4H-1-Benzopyran-4-one,3,7-dihydroxy-2-(3,4,5-trihydroxyphenyl)- CAS No.:490-31-3 Formula: C15H10O7 Molecular Weight : 302.25 Synonyms: Flavone,3,3',4',5',7-pentahydroxy- (8CI); Robinetin (6CI);3,3',4',5',7-Pentahydroxyflavone; 3,7,3',4',5'-Pentahydroxyflavone; NSC 407331;NSC 656274; Norkanugin EINECS:

Card-20(22)-enolide,3,14-dihydroxy-,(57263086,3b,5b)- CAS 143-62-4

Name: Card-20(22)-enolide,3,14-dihydroxy-,(57263087,3b,5b)- CAS No.:143-62-4 Formula: C23H34 O4 Molecular Weight : 374.57 Synonyms: 5b-Card-20(22)-enolide, 3b,14-dihydroxy- (6CI,7CI,8CI);(+)-Digitoxigenin; Cerberigenin; Digitoxigenin; NSC 407806; D20:22-3,14,21-Trihydroxynorcholenicacid lactone Density:

Ethylene Glycol CAS 107-21-1

Name: Ethylene glycol CAS No.:107-21-1 Formula: C2H6O2 Molecular Weight : 62.07 Synonyms: Glycol;Norkool;Ethylene dihydrate;1,2-Dihydroxyethane;Monoethylene glycol;146AR;M.e.g.;Ethylenglykol;Monoethylenglykol tech.;Ethyleneglycol;Glycol alcohol;Ethylene

L-Hydroxyproline CAS 51-35-4

Name: L-Hydroxyproline CAS No.:51-35-4 Formula: C5H9NO3 Molecular Weight : 131.13 Synonyms: Hydroxy-L-proline;(2S,4R)-4-hydroxypyrrolidine-2-carboxylic acid;trans-4-Hydroxyproline;L-Proline, 4-hydroxy-,

Propane,2-methoxy-2-methyl- CAS 1634-04-4

Name: Propane,2-methoxy-2-methyl- CAS No.:1634-04-4 Formula: C5H12O Molecular Weight : 88.15 Synonyms: Ether,tert-butyl methyl (6CI,7CI,8CI);1,1-Dimethylethyl methyl ether;2-Methoxy-2-methylpropane;2-Methyl-2-methoxypropane;MTBE;Methyl 1,1-dimethylethylether;Methyl tert-butyl ether;Methyl tertiary butyl ether;t-Butyl

2-Chloromethyl-3,4-dimethoxypyridinium chloride CAS 72830-09-2

Name: 2-Chloromethyl-3,4-dimethoxypyridinium chloride CAS No.:72830-09-2: Formula: C8H11Cl2NO2 Molecular Weight : 224.08 Synonyms: Pyridine,2-(chloromethyl)-3,4-dimethoxy-, hydrochloride (9CI); EINECS: 416-440-5 Boiling Point: 293.9 °C at 760 mmHg Flash Point: 131.6 °C Hazard Symbols: Xn, N

1,3:2,4-Bis(3,4-dimethylobenzylideno) sorbitol CAS 135861-56-2

Name: 1,3:2,4-Bis(3,4-dimethylobenzylideno) sorbitol CAS No.:135861-56-2 Formula: C24H30O6 Molecular Weight : 414.49 Synonyms: D-Glucitol,1,3:2,4-bis-O-[(3,4-dimethylphenyl) methylene]-;Millad 3988;Nucleating Agent NA-3;di-(3,4-dimethylbenzylidene)

N,O-Bis(trimethylsilyl)acetamide CAS 10416-59-8

Name: N,O-Bis(trimethylsilyl)acetamide CAS No.:10416-59-8 Formula: C8H21NOSi2 Molecular Weight : 203.43 Synonyms: Ethanimidic acid, N-(trimethylsilyl)-, trimethylsilyl ester;N-trimethylsilyl-1-trimethylsilyloxy-ethanimine;N-(Trimethylsilyl)acetimidic acid, trimethylsilyl ester;Trimethylsilyl

Lenalidomide CAS 191732-72-6

Name: Lenalidomide CAS No.:191732-72-6 Formula: C13H13N3O3 Molecular Weight : 259.26 Synonyms: Lenalidomide(other anti-cancers);Revlimid (lenalidomide);Revlimid (lenalidomide)/191732-72-6;Lenalidomide ; 3-(7-Amino-3-oxo-1H-isoindol-2-yl)piperidine-2,6-dione;Revlimid; Density: 1.46 g/cm3 Boiling Point: 614 °C at 760

Octreotide CAS 79517-01-4

Name: Octreotide CAS No.:79517-01-4 Formula: C49H66N10O10S2 Molecular Weight : 1019.24 Synonyms: L-Cysteinamide,D-phenylalanyl-L-cysteinyl-Lphenylalanyl- D-tryptophyl-L-lysyl-Lthreonyl- N-[(1R,2R)-2-hydroxy-1-(hydroxymethyl) propyl]-,cyclic (2f7)-disulfide,acetate (salt);Sandostatin (TN);Octreotide acetate

Olean-12-en-29-oicacid, 3-(3-carboxy-1-oxopropoxy)-11-oxo-,(57263075,3b,20b)- CAS 5697-56-3

Name: Olean-12-en-29-oicacid, 3-(3-carboxy-1-oxopropoxy)-11-oxo-,(57263076,3b,20b)- CAS No.:5697-56-3 Formula: C34H50O7 Molecular Weight : 570.76 Synonyms: Olean-12-en-30-oicacid, 3b-hydroxy-11-oxo-, hydrogensuccinate (7CI,8CI);Olean-12-en-30-oic acid, 3b-hydroxy-11-oxo-, succinate

1-(2-Hydroxyethyl)-2-imidazolidinone CAS 3699-54-5

Name: 1-(2-Hydroxyethyl)-2-imidazolidinone CAS No.:3699-54-5 Formula: C5H10N2O2 Molecular Weight : 130.14 Synonyms: 1-(2-Hydroxyethyl)-2-imidazolidinone;1-(Hydroxyethyl)ethyleneurea;N-(2-Hydroxyethyl)ethyleneurea;N-(2-Hydroxyethyl)imidazolidone;N-(b-Hydroxyethyl)-N,N'-ethylene urea; EINECS: 223-032-9 Density: 1.212

Benzenemethanol,3-bromo- CAS 15852-73-0

Name: Benzenemethanol,3-bromo- CAS No.:15852-73-0 Formula: C7H7BrO Molecular Weight : 187.03 Synonyms: Benzylalcohol, m-bromo- (6CI,8CI);3-Bromophenylmethanol;m-Bromobenzyl alcohol;M-Bromobenzyl Alcohol -1000Kgs; EINECS: 239-975-4 Density: 1.565 g/cm3 Boiling Point: 287.2 °C at 760 mmHg Flash Point: 115.7

DEXPANTHENOL CAS 81-13-0

Name: Dexpanthenol CAS No.:81-13-0 Formula: C9H19NO4 Molecular Weight : 205.25 Synonyms: Butanamide,2,4-dihydroxy-N-(3-hydroxypropyl)-3,3-dimethyl-,(57263069,R)-;Butyramide,2,4-dihydroxy-N-(3-hydroxypropyl)-3,3-dimethyl-, D-(+)- (8CI);(+)-Panthenol;Alcopan

Chromium carbonyl(Cr(CO)6),(57263066,OC-6-11)- CAS 13007-92-6

Name: Chromium carbonyl(Cr(CO)6),(57263067,OC-6-11)- CAS No.:13007-92-6 Formula: C6CrO6 Molecular Weight : 220.06 Synonyms: Chromiumcarbonyl (Cr(CO)6) (8CI);Chromium carbonyl;Hexacarbonylchromium; EINECS: 235-852-4 Density: 1.77 g/cm3 Boiling Point: 220 °C Solubility: Insoluble in water Appearance: white crystals or

Lactitol Monohydrate CAS 81025-04-9

Name: D-Glucitol, 4-O-b-D-galactopyranosyl-, hydrate(1:1) CAS No.:81025-04-9 Formula: C12H26O12 Molecular Weight : 362.38 Synonyms: D-Glucitol,4-O-b-D-galactopyranosyl-, monohydrate(9CI); EINECS: 209-566-5 Density: 1.69 g/cm3 Boiling Point: 788.5 °C at 760 mmHg Flash Point: 430.7 °C Solubility: Soluble in

Voglibose CAS 83480-29-9

Name: Voglibose CAS No.:83480-29-9 Formula: C10H21NO7 Molecular Weight : 267.28 Synonyms: Vioglibose;n-(1,3-dihydroxy-2-propyl)valiolamine;5-(1,3-dihydroxypropan-2-ylamino)-1-(hydroxymethyl)cyclohexane-1,2,3,4-tetrol;AO 128;Voglibose (JAN/USAN);Basen (TN);A

[1,1'-Biphenyl]-4-aceticacid, 2-fluoro-a-methyl-,1-(acetyloxy)ethyl ester CAS 91503-79-6

Name: [1,1'-Biphenyl]-4-aceticacid, 2-fluoro-a-methyl-,1-(acetyloxy)ethyl ester CAS No.:91503-79-6 Formula: C19H19 F O4 Molecular Weight : 330.38 Synonyms: 1-Acetoxyethyl2-(2-fluoro-4-biphenylyl)propionate;Flurbiprofen axetil;LFP 83;Ropiopn;1-Acetoxyethyl 2-(2-fluorobiphenyl-4-yl)propanoate; Density:

2-Acetylpyridine CAS 1122-62-9

Name: 2-Acetylpyridine CAS No.:1122-62-9 Formula: C7H7NO Molecular Weight : 121.15 Synonyms: FEMA 3251;2-Acetopyridine;1-(2-Pyridinyl)ethanone;Ethanone, 1-(2-pyridinyl)-;1-pyridin-2-ylethanone;Methyl 2-pyridyl ketone;Ketone, methyl 2-pyridyl;2-Acetylpyridine (natural);2-Pyridyl methyl ketone;2-Acetyl pyridine;Acetyl

5,6-Dihydro-4-hydroxy-6-methyl-4H-thieno[2,3-b]thiopyran-2-sulfonamide 7,7-dioxide CAS 120279-26-7

Name: 5,6-Dihydro-4-hydroxy-6-methyl-4H-thieno[2,3-b]thiopyran-2-sulfonamide 7,7-dioxide CAS No.:120279-26-7 Formula: C8H11NO5S3 Molecular Weight : 297.37 Synonyms: INTERMEDIATE FOR DORZOLAMIDE;5,6-DIHYDRO-4H-4-HYDROXY-6-METHYL THIENO [2,3,B]

Thiram CAS 137-26-8

Name: Thiram CAS No.:137-26-8 Formula: C6H12N2S4 Molecular Weight : 240.432880 g/mol Synonyms: Rubber Accelerator;Thioperoxydicarbonicdiamide ([(H2N)C(S)]2S2), N,N,N',N'-tetramethyl-;Disulfide,bis(dimethylthiocarbamoyl) (8CI);AApirol;Aatiram;Accel TMT;Accel TT;Accelerator Thiuram;Anles;Arasan 50 red;Arasan 70-S

Darifenacin hydrobromide CAS 133099-07-7

Name: Darifenacin hydrobromide CAS No.:133099-07-7 Formula: C28H30N2O2.HBr Molecular Weight : 507.47 Synonyms: 2-[(3S)-1-[2-(2,3-dihydrobenzofuran-5-yl)ethyl]pyrrolidin-3-yl]-2,2-diphenyl-acetamide hydrobromide;Enablex (TN);3-Pyrrolidineacetamide, 1-(2-(2,3-dihydro-5-benzofuranyl)ethyl)-alpha,alpha-diphenyl-,

2-Propenamide,2-cyano-3-(3,4-dihydroxy-5-nitrophenyl)-N,N-diethyl-,(57263054,2E)- CAS 130929-57-6

Name: 2-Propenamide,2-cyano-3-(3,4-dihydroxy-5-nitrophenyl)-N,N-diethyl-,(57263055,2E)- CAS No.:130929-57-6 Formula: C14H15N3O5 Molecular Weight : 305.29 Synonyms: Entacom;OR 611;2-Propenamide,2-cyano-3-(3,4-dihydroxy-5-nitrophenyl)-N,N-diethyl-,(57263056,E)-;(E)-Entacapone;Comtan; Density: 1.392 g/cm3 Boiling

Cilostazol CAS 73963-72-1

Name: Cilostazol CAS No.:73963-72-1 Formula: C20H27N5O2 Molecular Weight : 369.52 Synonyms: 6-[4-(1-cyclohexyltetrazol-5-yl)butoxy]-3,4-dihydro-1H-quinolin-2-one;Pletal;OPC 13013;Cilostazol (JAN/USAN);2(1H)-Quinolinone,6-[4-(1-cyclohexyl-1Htetrazol- 5-yl)butoxy]-3,4-dihydro-;Pletal (TN);Cilostazole;Cilostal; Density:

Propanoic acid,3-(benzoylthio)-2-methyl- CAS 74431-50-8

Name: Propanoic acid,3-(benzoylthio)-2-methyl- CAS No.:74431-50-8 Formula: C11H12O3S Molecular Weight : 224.28 Synonyms: Propanoic acid,3-(benzoylthio)-2-methyl-,(57263051,?à)-;3-(Benzoylthio)-2-methylpropionic acid;3-(Benzoylthio)isobutyric acid; EINECS: 277-867-9 Density: 1.249 g/cm3 Boiling Point: 367.5 °C at 760

4-Bromoindole CAS 52488-36-5

Name: 4-Bromoindole CAS No.:52488-36-5 Formula: C8H6BrN Molecular Weight : 196.05 Synonyms: 4-bromo-1H-indole; EINECS: -0 Density: 1.66 g/cm3 Boiling Point: 316.9 °C at 760 mmHg Flash Point: 145.5 °C Appearance: Clear yellowish-green or dark brown liquid

ZINC SALICYLATE CAS 16283-36-6

Name: Zinc, bis[2-(hydroxy-kO)benzoato-kO]-,(57263047,T-4)- CAS No.:16283-36-6 Formula: Zn.(C7H5O3)2 Molecular Weight : 339.61 Synonyms: Zinc,bis(2-hydroxybenzoato-O1,O2)-,(57263048,T-4)-;Zinc, bis(salicylato)- (6CI,7CI,8CI);Bis(salicylato)zinc;Bontron E 304;DNA (Fugu rubripes clone 024P08bG3 genomesurvey

6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol,4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-,(57263044,4aS,6R,8aS)- CAS 357-70-0

Name: 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol,4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-,(57263045,4aS,6R,8aS)- CAS No.:357-70-0 Formula: C17H21NO3 Molecular Weight : 287.35 Synonyms: 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol,4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl- (7CI);Galanthamine

3(2H)-Pyridazinone,4,5-dihydro-6-[2-(4-methoxyphenyl)-1H-benzimidazol-6-yl]-5-methyl- CAS 74150-27-9

Name: 3(2H)-Pyridazinone,4,5-dihydro-6-[2-(4-methoxyphenyl)-1H-benzimidazol-6-yl]-5-methyl- CAS No.:74150-27-9 Formula: C19H18N4O2 Molecular Weight : 334.37 Synonyms: Acardi;Racemic pimobendan;UD-CG 115;UD-CG 115BS;Vetmedin;3(2H)-Pyridazinone,4,5-dihydro-6-[2-(4-methoxyphenyl)-1H-benzimidazol-5-yl]-5-methyl- (9CI);

Ethylenediaminetetraacetic acid CAS 60-00-4

Name: Ethylenediaminetetraacetic acid CAS No.:60-00-4 Formula: C10H16N2O8 Molecular Weight : 292.28 Synonyms: Glycine,N,N'-1,2-ethanediylbis[N-(carboxymethyl)-;Aceticacid,(57263041,ethylenedinitrilo)tetra-(8CI);3,6-Diazaoctanedioic acid,3,6-bis(carboxymethyl)-;Cheelox;Complexon II;EDTA (chelating

C.I. Acid Blue 82 CAS 12217-19-5

Name: C.I. Acid Blue 82 CAS No.:12217-19-5 Formula: Unspecified Molecular Weight : 0 Synonyms: Acid Blue82;Acid Fast Light Blue JNL;Best Acid Blue GW;Coomassie Blue P;Eriosin FastBlue GL;Mitsui Nylon Fast Blue B;Sulphonol Fast BlueP;Suminol Fast Blue G;Superlan Night Blue AG;Supramin Blue GW;TertracidFast Blue

Benzoic acid,3-(aminosulfonyl)-4-chloro-5-nitro- CAS 22892-96-2

Name: Benzoic acid,3-(aminosulfonyl)-4-chloro-5-nitro- CAS No.:22892-96-2 Formula: C7H5ClN2O6S Molecular Weight : 280.64 Synonyms: Benzoicacid, 4-chloro-3-nitro-5-sulfamoyl- (8CI);4-Chloro-3-nitro-5-sulfamoylbenzoateacid;4-Chloro-3-nitro-5-sulfamoylbenzoic acid;4-Chloro-3-nitro-5-sulfamylbenzoic

D-Glucose, 6-O-a-D-glucopyranosyl- CAS 499-40-1

Name: D-Glucose, 6-O-a-D-glucopyranosyl- CAS No.:499-40-1 Formula: C12H22O11 Molecular Weight : 342.2965 Synonyms: Isomaltose(8CI);6-O-a-D-Glucopyranosyl-D-glucose;D-Isomaltose; EINECS: 207-879-1 Density: 1.68 g/cm3 Boiling Point: 774.5 °C at 760 mmHg Flash Point: 288.8 °C

Linoleic acid CAS 60-33-3

Name: 9,12-Octadecadienoicacid (9Z,12Z)- CAS No.:60-33-3 Formula: C18H32O2 Molecular Weight : 280.50 Synonyms: 9,12-Octadecadienoicacid (Z,Z)-;Linoleic acid (8CI);(9Z,12Z)-9,12-Octadecadienoic acid;(Z,Z)-9,12-Octadecadienoic acid;9,12-Octadecadienoic acid,(57263036,Z,Z)-;9-cis,12-cis-Linoleic acid;9Z,12Z-Linoleic

Pyridine,2,3-dichloro-5-(trichloromethyl)- CAS 69045-83-6

Name: Pyridine,2,3-dichloro-5-(trichloromethyl)- CAS No.:69045-83-6 Formula: C6H2Cl5N Molecular Weight : 265.35 Synonyms: 2,3-Dichloro-5-trichloromethylpyridine; Density: 1.681 g/cm3 Boiling Point: 293.229 °C at 760 mmHg Flash Point: 158.979 °C

Ferrous lactate CAS 5905-52-2

Name: Ferrous lactate CAS No.:5905-52-2 Formula: C6H10FeO6 Molecular Weight : 234.01 Synonyms: Lacticacid, iron(2+) salt (2:1) (8CI);CromatonbicFerro;E 585;Ferro-Drops;Ferrouslactate; EINECS: 227-608-0 Boiling Point: 227.6 °C at 760 mmHg Flash Point: 109.9 °C Appearance: green-white powder

2-Methyl-octahydro-pyrrolo(3,4-c)pyridine CAS 16039-64-8

Name: 2-Methyl-octahydro-pyrrolo(3,4-c)pyridine CAS No.:16039-64-8 Molecular Weight : 457.00000 Synonyms: potassium,antimony(3+),2,3-dihydroxybutanedioate;Potassium Antimony Tartrate; Boiling Point: 399.3oC at 760 mmHg Flash Point: 209.4oC

Riboflavin CAS 83-88-5

Name: Riboflavin CAS No.:83-88-5 Formula: C17H20N4O6 Molecular Weight : 376.37 . Synonyms: Vitamin B2-Riboflavin(Vitamin B2)USP/BP/EP;VB2

Potassium nitrate CAS 7757-79-1

Name: Potassium nitrate CAS No.:7757-79-1 Formula: KNO3 Molecular Weight : 101.10 Synonyms: Nitric acid, potassium salt;Nitric acid potassium salt;Potassium Nitrate ACS Crystals;Nitre;Salt

peter;Sensodyne;Niter;Sensodyne (TN);Potassium nitrate (JAN/USP);Collo-Bo;Saltpeter;Potassium nitrate(Pharmaceutical

Phosphonic acid,[[(phosphonomethyl)imino]bis[2,1-ethanediylnitrilobis(methylene)]]tetrakis-,sodium salt (1:?) CAS 22042-96-2

Name: Phosphonic acid,[[(phosphonomethyl)imino]bis[2,1-ethanediylnitrilobis(methylene)]]tetrakis-,sodium salt (1:?) CAS No.:22042-96-2 Formula: C9H28N3O15P5.xNa Molecular Weight : 793.01700 Synonyms: Phosphonic acid,[[bis[2-[bis(phosphonomethyl)amino]ethyl]amino]methyl]-, sodium salt (8CI);Briquest 221;Briquest

3-Aminopyridine-4-carboxylic acid methyl ester CAS 55279-30-6

Name: 3-Aminopyridine-4-carboxylic acid methyl ester CAS No.:55279-30-6 Formula: C7H8N2O2 Molecular Weight : 152.15 Synonyms: 3-Aminoisonicotinicacid methyl ester;Methyl3-amino-4-pyridinecarboxylate;Methyl 3-aminoisonicotinate;4-Pyridinecarboxylicacid, 3-amino-, methyl ester; Density: 1.238 g/cm3 Boiling Point: 292.5

Butanoic acid,1,1'-anhydride CAS 106-31-0

Name: Butanoic acid,1,1'-anhydride CAS No.:106-31-0 Formula: C8H14O3 Molecular Weight : 158.19 Synonyms: Butanoicacid, anhydride (9CI);Butyric anhydride (6CI,8CI);Butanoic anhydride;Butanoyl anhydride;Butyric acid anhydride;Butyryl oxide;n-Butyric acidanhydride;n-Butyric anhydride; EINECS: 203-383-4 Density: 0.981

Lanosta-8,24-dien-3-ol,(57263025,3b)- CAS 79-63-0

Name: Lanosta-8,24-dien-3-ol,(57263026,3b)- CAS No.:79-63-0 Formula: C30H50O Molecular Weight : 426.72 Synonyms: Lanosta-8,24-dien-3b-ol (8CI);Lanostadien-3b-ol (7CI);Lanosterol

Flupirtine CAS 56995-20-1

Name: Carbamic acid,N-[2-amino-6-[[(4-fluorophenyl)methyl]amino]-3-pyridinyl]-, ethyl ester CAS No.:56995-20-1 Formula: C15H17FN4O2 Molecular Weight : 304.32 Synonyms: Carbamicacid, [2-amino-6-[[(4-fluorophenyl)methyl]amino]-3-pyridinyl]-, ethyl ester(9CI);D 9998;Flupirtine;Katadolon;Trancopal Dolo; EINECS:

Deoxyribonuclease CAS 9003-98-9

Name: Deoxyribonuclease CAS No.:9003-98-9 Formula: Molecular Weight : 0 Synonyms: AlkalineDNase;Alkaline deoxyribonuclease;DNA depolymerase;DNA endonuclease;DNAnuclease;DNAase;DNase;DNase I;DNase Y;DNase g;Deoxyribonuclease;Deoxyribonuclease (pancreatic);Deoxyribonuclease A;Deoxyribonucleic

1,3,5-Triazine-2,4,6(1H,3H,5H)-trione, 1,3,5-tris((3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl)methyl)- CAS 27676-62-6

Name: 1,3,5-Triazine-2,4,6(1H,3H,5H)-trione, 1,3,5-tris((3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl)methyl)- CAS No.:27676-62-6 Formula: C48H69N3O6 Molecular Weight : 784.08 Synonyms: s-Triazine-2,4,6(1H,3H,5H)-trione,1,3,5-tris(3,5-di-tert-butyl-4-hydroxybenzyl)- (8CI);1,3,5-Tris(3,5-di-tert-butyl-4-hydroxybenzyl)

1H-Pyrazolo[3,4-d]pyrimidin-4-amine,3-iodo- CAS 151266-23-8

Name: 1H-Pyrazolo[3,4-d]pyrimidin-4-amine,3-iodo- CAS No.:151266-23-8 Formula: C5H4IN5 Molecular Weight : 261.02 Synonyms: 3-iodo-4-amino-1H-pyrazolo[3,4-d]pyrimidine;3-Iodo-1H-pyrazolo[3,4-d]pyrimidin-4-amine; Density: 2.465 g/cm3 Boiling Point: 498.8 °C at 760 mmHg Flash Point: 255.5 °C

XYLENE CAS 1330-20-7

Name: Xylene CAS No.:1330-20-7 Formula: C8H10 Molecular Weight : 106.18 Synonyms: Dimethylbenzene;Xylene mixture (60% m-xylene, 9% o-xylene, 14% p-xylene, 17% ethylbenzene);Xylenes (mixed);Xylol;Benzene,dimethyl-;Eylene; EINECS: 215-535-7 Density: 0.865 g/cm3 Boiling Point: 136-140 ºC Flash Point: 21 ºC Solubility:

Propyne CAS 74-99-7

Name: Propyne CAS No.:74-99-7 Formula: C3H4 Molecular Weight : 40.07 Synonyms: Propyne(8CI);Allylene;Methylacetylene;1-Propyne; EINECS: 200-828-4 Density: 0.648 g/cm3 Boiling Point: -23.2 °C(lit.) Flash Point: -51 °C

3,4-Dihydroxybenzonitrile CAS 17345-61-8

Name: 3,4-Dihydroxybenzonitrile CAS No.:17345-61-8 Formula: C7H5NO2 Molecular Weight : 135.12 Synonyms: 3,4-Dihydroxyl benzonitrile 98%;3,4-Dihydroxy benzonitrile;4-Cyanocatechol; EINECS: 241-367-9

Density: 1.42 g/cm3 Boiling Point: 334.8 °C at 760 mmHg Flash Point: 156.3 °C Solubility: insoluble in water Appearance:

Lincomycin hydrochloride CAS 859-18-7

Name: Lincomycin hydrochloride CAS No.:859-18-7 Formula: C18H34N2O6S.HCl Molecular Weight : 443.06 Synonyms: D-erythro-D-galacto-Octopyranoside,methyl 6,8-dideoxy-6-(1-methyl-4-propyl-L-2-pyrrolidinecarboxamido)-1-thio-,monohydrochloride, trans-a- (8CI);D-erythro-a-D-galacto-Octopyranoside,

Methyl 4-methylbenzoate CAS 99-75-2

Name: Methyl 4-methylbenzoate CAS No.:99-75-2 Formula: C9H10O2 Molecular Weight : 150.19 Synonyms: p-Toluicacid, methyl ester (6CI,8CI);4-(Methoxycarbonyl)toluene;4-Methylbenzoic acidmethyl ester; EINECS: 202-784-1 Density: 1.045 g/cm3 Boiling Point: 222.5 °C at 760 mmHg Flash Point: 90.6 °C Solubility:

7-Keto-dehydroepiandrosterone CAS 566-19-8

Name: 7-Keto-dehydroepiandrosterone CAS No.:566-19-8 Formula: C19H26O3 Molecular Weight : 304.43 Synonyms: Androst-5-ene-7,17-dione,3b-hydroxy- (8CI);3b-Hydroxy-5-androstene-7,17-dione;5-Androsten-3b-ol-7,17-dione;7-Keto-DHEA;7-Oxo-DHEA;Androst-5-ene-7,17-dione,3-hydroxy-,(57263408,3b)-; Density: 1.19 g/cm3 Boiling

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Россия +7(495)268-04-70 Казахстан +7(7172)727-132 Киргизия +996(312)96-26-47

эл.почта: tuf@nt-rt.ru || сайт: https://tianfu.nt-rt.ru