

# Audit - EU DK MAL Code

## PPG AQUACOVER 45 NF BASE L

### Denmark MAL Code

#### Audit - MAL Code

EU Denmark MAL Code:- 00-3

The MAL Code calculations are performed with product and component data.

Product is a Liquid

PPG AQUACOVER 45 NF BASE L - Components considered for the MAL Code calculation. {Denmark MAL Code}

WATER (46.283521446344%)

CAS: 7732-18-5

Density: 1

Molecular Weight: 18.02

Boiling Point: 100

Vapour Pressure: 17.5

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 0; Lower Limit: 0

proprietary acrylic copolymer (26.32282%)

CAS: SUB122235

Density: 0

No LBL Factor entered or estimated from CAS Number or Boiling Point.

No MAL Factor calculated.

FAD: 1. (Default)

FAD 1 Quotient = 26322.82

TITANIUM DIOXIDE (18.0343695773922%)

CAS: 13463-67-7

Density: 4.1

Relative Density: 4.26

Molecular Weight: 79.9

Boiling Point: 2750

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: 0

FAD 1 Quotient = 18034.370

2-(2-BUTOXYETHOXY)ETHANOL (1.8632053%)

CAS: 112-34-5

Density: 0.953

Relative Density: 0.95

Molecular Weight: 162.26

Boiling Point: 226.3

Vapour Pressure: 0.02

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.186

SODIUM POTASSIUM ALUMINUM SILICATE (1.5%)

CAS: 37244-96-5

Density: 2.56

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: 0.1

FAD 1 Quotient = 15

POLYPROPYLENE GLYCOL (1.4525%)

CAS: 25322-69-4

Density: 0

Relative Density: 1.01

Vapour Pressure: 0.00063

No LBL Factor entered or estimated from CAS Number or Boiling Point.

R Phrases: Xn;R22

MAL Factor from Sub-Annex 2: 0

FAD: 3. (Xn)

FAD 3 Quotient = 1452.5

ALUMINUM HYDROXIDE (0.76002275%)

CAS: 21645-51-2

Density: 2.42

Molecular Weight: 78

Vapour Pressure: 0.0675

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: 0.1

FAD 1 Quotient = 7.600

POLYMER, POLYFUNCTIONAL, NON-ANIONIC (0.4419983%)

CAS: SUB110823

Density: 1.1

No LBL Factor entered or estimated from CAS Number or Boiling Point.

No MAL Factor calculated.

FAD: 1. (Default)

FAD 1 Quotient = 441.998

POLYURETHANE RESIN (0.394235%)

CAS: SUB100112

Density: 1.1

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: 0.1

FAD 1 Quotient = 3.942

COALESCING AID (0.3496479%)

CAS: SUB106738

Density: 0

No LBL Factor entered or estimated from CAS Number or Boiling Point.

No MAL Factor calculated.

FAD: 1. (Default)

FAD 1 Quotient = 349.648  
Alcohols, C16-18 and C18-unsatd., ethoxylated (0.34023078279291%)  
CAS: 68920-66-1  
Density: 1  
Boiling Point: 369  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: 0.1  
FAD 1 Quotient = 3.402  
POLYSILOXANE MIXTURE (0.297362275%)  
CAS: SUB100136  
Density: 1.1  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 297.362  
Tripropylene glycol monomethyl ether (0.29155%)  
CAS: 25498-49-1  
Density: 0.96  
Molecular Weight: 206.32  
Boiling Point: 242.8  
Vapour Pressure: 0.01500123  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 291.55  
proprietary surfactant (0.28003%)  
CAS: SUB122236  
Density: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 280.03  
POLYACRYLATE (0.2298666%)  
CAS: SUB117312  
Density: 1.09  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 229.867  
Methanone, (1-hydroxycyclohexyl)phenyl- (0.2%)  
CAS: 947-19-3  
Density: 1.17  
Boiling Point: 316.1  
Vapour Pressure: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: 0.1

FAD 1 Quotient = 2  
BENZOPHENONE (0.1996%)  
CAS: 119-61-9  
Density: 1.09  
Relative Density: 1.1  
Molecular Weight: 182.23  
Boiling Point: 305.4  
Vapour Pressure: 0.003  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.100

DIPROPYLENE GLYCOL MONOMETHYL ETHER (0.11645%)  
Organic Solvent.  
CAS: 34590-94-8  
Density: 0.95  
Relative Density: 0.95  
Molecular Weight: 148.23  
Boiling Point: 189.6  
Vapour Pressure: 0.277522755  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 5. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 116.45

TRIMETHYLOLPROPANE (0.085312925%)  
CAS: 77-99-6  
Density: 1.084  
Molecular Weight: 134.2  
Boiling Point: 304.2  
Vapour Pressure: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.853

SODIUM NITRITE (0.064%)  
CAS: 7632-00-0  
Density: 2.2  
Relative Density: 2.17  
Molecular Weight: 69  
Boiling Point: 320  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.32  
FAD 3 Quotient = 0.64

3-Iodo-2-propynyl butylcarbamate (0.0599421590592%)  
CAS: 55406-53-6  
Density: 1.5

Molecular Weight: 281.11  
Vapour Pressure: 0.0000063  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.060

ZIRCONIUM TETRAHYDROXIDE (0.057%)

CAS: 14475-63-9  
Density: 1.5  
Molecular Weight: 159.25  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.57

reaction mass of mixed (3,3,4,4,5,5,6,6,7,7, 8,8,8- tridecafluorooctyl) phosphates, ammonium salt (0.0544%)

CAS: SUB141402  
Density: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 54.4

AMMONIUM HYDROXIDE (0.04571%)

CAS: 1336-21-6  
Density: 0.9  
Relative Density: 0.9  
Molecular Weight: 35.06  
Boiling Point: 38  
Vapour Pressure: 360.02925  
LB�Factor = 100 (BP=38)  
MAL Factor entered: 50. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 4 Quotient = 0.001  
FAD 3 Quotient = 0.009

polyurethane resin (0.0425%)

CAS: SUB142197  
Density: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 42.5

AMMONIUM BENZOATE (0.04%)

CAS: 1863-63-4  
Density: 1.26  
Relative Density: 1.26  
Molecular Weight: 139.15  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.04  
proprietary defoamer (0.0392042%)  
CAS: SUB122237  
Density: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 39.204  
POLYETHYLENE-POLYPROPYLENE POLYMER (0.03780342031032%)  
CAS: 9003-11-6  
Density: 1.1  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.378  
1,2-BENZISOTHAZOLONE (0.0330321585497%)  
CAS: 2634-33-5  
Density: 1.095  
Molecular Weight: 151.19  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.033  
DIMETHYLAMINOETHANOL (0.028%)  
Organic Solvent.  
CAS: 108-01-0  
Density: 0.89  
Relative Density: 0.89  
Molecular Weight: 89.14  
Boiling Point: 134.1  
Vapour Pressure: 4.59  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 280. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.003  
FAD 2 Quotient = 0.014  
TITANIUM DIOXIDE (<10 microns) (0.016245%)  
Carcinogen.  
CAS: 13463-67-7  
Density: 4.1  
Relative Density: 4.26  
Molecular Weight: 79.9  
Boiling Point: 2750  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 16.245  
ZINC OXIDE (0.015%)

CAS: 1314-13-2

Density: 5.68

Relative Density: 5.61

Molecular Weight: 81.37

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: 0

FAD 1 Quotient = 15

polyethylene glycol monobutyl ether (0.00925%)

CAS: 90736-95-1

Density: 0

No LBL Factor entered or estimated from CAS Number or Boiling Point.

No MAL Factor calculated.

FAD: 1. (Default)

FAD 1 Quotient = 9.25

2-BUTOXY ETHANOL (0.0055823%)

Organic Solvent.

CAS: 111-76-2

Density: 0.9

Relative Density: 0.9

Molecular Weight: 118.18

Boiling Point: 171.25

Vapour Pressure: 0.75006

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 25. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.001

pyrithione zinc (0.0055%)

CAS: 13463-41-7

Density: 1.76

Molecular Weight: 317.69

Boiling Point: 269.85

Vapour Pressure: 0.0000000072

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.006

1-BUTANOL (0.00185%)

Organic Solvent.

CAS: 71-36-3

Density: 0.81

Relative Density: 0.81

Molecular Weight: 74.14

Boiling Point: 119

Vapour Pressure: 6.750576

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 67. Limit: 0

FAD entered: 1; Lower Limit: 0

FAD 1 Quotient = 1.85  
ISOTHIAZOLONE SOLUTION (0.00086624186471%)  
CAS: 55965-84-9  
Density: 0.9  
Molecular Weight: 264.76  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 0.866  
HEPTANE (0.000396%)  
Organic Solvent.  
CAS: 142-82-5  
Density: 0.684  
Relative Density: 0.68  
Molecular Weight: 100.23  
Boiling Point: 98.5  
Vapour Pressure: 34.5028  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 12. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 0.396  
DIETHYLENE GLYCOL (0.00019057%)  
Organic Solvent.  
CAS: 111-46-6  
Density: 1.18  
Relative Density: 1.12  
Molecular Weight: 106.12  
Boiling Point: 244.9  
Vapour Pressure: 0.006  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.000  
ACETIC ACID (0.000188232049%)  
Organic Solvent.  
CAS: 64-19-7  
Density: 1.04  
Relative Density: 1.05  
Molecular Weight: 60.06  
Boiling Point: 117.9  
Vapour Pressure: 15.59383  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 1. Limit:  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 4 Quotient = 0.000  
Triethyleneglycol monobutylether (0.000153%)  
CAS: 143-22-6  
Density: 0.99



Relative Density: 0.98  
Molecular Weight: 206.32  
Boiling Point: 278  
Vapour Pressure: 0.0075  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.002  
FAD 3 Quotient = 0.000

residual monomers (0.0001334%)

CAS: SUB137626  
Density: 0  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 0.133

PROPYLENE OXIDE (0.0000525%)

Organic Solvent.  
Carcinogen.  
CAS: 75-56-9  
Density: 0.83  
Relative Density: 0.8  
Molecular Weight: 58.09  
Boiling Point: 34.23  
Vapour Pressure: 538  
LBLFactor = 100 (BP=34.23)  
MAL Factor entered: 1. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.000

TRIETHYLENEGLYCOL (0.00004089609079%)

CAS: 112-27-6  
Density: 1.125  
Relative Density: 1.1  
Molecular Weight: 150.2  
Boiling Point: 286.5  
Vapour Pressure: 0.00049  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.000

2-METHYL-4-ISOTHIAZOLIN-3-ONE (0.00003588009408%)

CAS: 2682-20-4  
Density: 0.8  
Molecular Weight: 115.1  
Boiling Point: 94  
Vapour Pressure: 0.000037503  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 6 Quotient = 0.000

FAD 3 Quotient = 0.001

OCTAMETHYLCYCLOTETRAILOXANE (0.0000280019274%)

CAS: 556-67-2

Density: 0.95

Relative Density: 0.96

Molecular Weight: 296.68

Boiling Point: 175

Vapour Pressure: 0.99008

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 1. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

Decamethylcyclopentasiloxane (0.0000280019274%)

CAS: 541-02-6

Density: 0.96

Molecular Weight: 370.85

Boiling Point: 210

Vapour Pressure: 0.25

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

dodecamethylcyclohexasiloxane (0.0000280019274%)

CAS: 540-97-6

Density: 0.98

Molecular Weight: 445.02

No LBL Factor entered or estimated from CAS Number or Boiling Point.

No MAL Factor calculated.

FAD: 1. (Default)

FAD 1 Quotient = 0.028

alkyl polyglycol ether phosphate compound (0.00002467448481%)

CAS: 164383-18-0

Density: 1.1

Boiling Point: 220

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

ETHANOL;2-(2-ETHOXYETHOXY) (0.0000153%)

CAS: 111-90-0

Density: 0.986

Relative Density: 0.99

Molecular Weight: 134.18

Boiling Point: 196

Vapour Pressure: 0.14

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

Triethylene glycol monoethyl ether (0.0000153%)

CAS: 112-50-5

Density: 1.021

Relative Density: 1.02

Molecular Weight: 178.26

Boiling Point: 256

Vapour Pressure: 0.01

No LBL Factor entered or estimated from CAS Number or Boiling Point.

R Phrases: None

MAL Factor from Sub-Annex 2: 0

FAD: 1. (Default)

FAD 1 Quotient = 0.015

ALUMINUM SILICATE (0.00001458122715%)

CAS: 1332-58-7

Density: 2.6

Relative Density: 2.6

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

AMORPHOUS SILICA (0.00001195212918%)

CAS: 112945-52-5

Density: 1.5

Molecular Weight: 60.09

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

2-ETHYLHEXANOIC ACID (0.00000735%)

CAS: 149-57-5

Density: 0.9

Relative Density: 0.9

Molecular Weight: 144.24

Boiling Point: 227.5

Vapour Pressure: 0.03

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

SILICA (0.00000650023922%)

CAS: 7631-86-9

Density: 2

Relative Density: 2.2

Molecular Weight: 60.08

Boiling Point: 2230

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

R Phrases: None

FAD: 1. (Default)

FAD 1 Quotient = 0.007

LECITHINS (0.00000528808926%)

CAS: 8002-43-5

Density: 1.1

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

Ethanol, 2,2'-(butylimino)bis- (0.00000444885%)

CAS: 102-79-4

Density: 0.968

Relative Density: 0.99

Molecular Weight: 161.28

Boiling Point: 274

Vapour Pressure: 0.877571955

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 1. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

BENZENE (0.000004%)

Organic Solvent.

Carcinogen.

CAS: 71-43-2

Density: 0.877

Relative Density: 0.88

Molecular Weight: 78.12

Boiling Point: 80.09

Vapour Pressure: 75.00609

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 880. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 6 Quotient = 0.000

ZIRCONIUM OXIDE (0.00000325%)

CAS: 1314-23-4

Density: 5.85

Molecular Weight: 123.22

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

TIN (0.0000017%)

CAS: 7440-31-5

Density: 7.2

Relative Density: 7.28

Molecular Weight: 118.69  
Boiling Point: 2260  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor from OEL: 0  
R Phrases: None  
FAD: 1. (Default)  
FAD 1 Quotient = 0.002

POLYETHER SILOXANE COPOLYMER (0.00000042488445%)

CAS: SUB117132  
Density: 1.1  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 0.000

SODIUM HYDROXIDE (0.0000004019274%)

CAS: 1310-73-2  
Density: 2.1  
Relative Density: 2.13  
Molecular Weight: 40  
Boiling Point: 1390  
Vapour Pressure: 0.097507995  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 4 Quotient = 0.000  
FAD 3 Quotient = 0.000

ACETALDEHYDE (0.00000035%)

Organic Solvent.  
Carcinogen.  
CAS: 75-07-0  
Density: 0  
Relative Density: 0.78  
Molecular Weight: 44.06  
Boiling Point: 20.1  
Vapour Pressure: 900.07313  
LBLFactor = 100 (BP=20.1)  
MAL Factor entered: 1. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.000

polycarbonic acid ammonium salt (0.00000030709786%)

CAS: SUB109712  
Density: 1.32  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
No MAL Factor calculated.  
FAD: 1. (Default)  
FAD 1 Quotient = 0.000

2-BROMO-2-NITRO-1,3-PROPANEDIOL (0.0000001988%)

CAS: 52-51-7

Density: 1.1  
Relative Density: 1.1  
Molecular Weight: 200.01  
Vapour Pressure: 0.00004  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.000

1,4-DIOXANE (0.000000175%)

Organic Solvent.  
Carcinogen.  
CAS: 123-91-1  
Density: 1.03  
Relative Density: 1.03  
Molecular Weight: 88.12  
Boiling Point: 101.15  
Vapour Pressure: 30.7525  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 390. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.000  
FAD 3 Quotient = 0.000

FORMALDEHYDE (0.000000175%)

Carcinogen.  
CAS: 50-00-0  
Density: 1.09  
Relative Density: 0.812  
Molecular Weight: 30.03  
Boiling Point: 98  
Vapour Pressure: 1  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 2500. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.000  
FAD 3 Quotient = 0.000

METHYL ALCOHOL (0.000000175%)

Organic Solvent.  
CAS: 67-56-1  
Density: 0.792  
Relative Density: 0.79  
Molecular Weight: 32.05  
Boiling Point: 64.7  
Vapour Pressure: 126.96329  
LBLFactor = 100 (BP=64.7)  
MAL Factor entered: 54. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.000  
FAD 3 Quotient = 0.000

ETHYLENE OXIDE (0.000000175%)

Carcinogen.

CAS: 75-21-8

Density: 0.882

Relative Density: 0.9

Molecular Weight: 44.06

Boiling Point: 10.7

Vapour Pressure: 1314.1117

LBLFactor = 100 (BP=10.7)

MAL Factor entered: 11. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 6 Quotient = 0.000

METHYL CHLORIDE (0.000000175%)

Carcinogen.

CAS: 74-87-3

Density: 0.911

Relative Density: 0.92

Molecular Weight: 50.49

Boiling Point: -23.7

Vapour Pressure: 3671.9

LBLFactor = 100 (BP=-23.7)

MAL Factor from OEL: 476.19 \*\* Warning: An Evaporation Rate Correction Factor of 2 was used. Contact the Authorities for a MAL Factor.

R Phrases: F+;R12 Xn;R48/20 Carc.Cat.3;R40

FAD: 1. (Default)

FAD 1 Quotient = 0.000

QUARTZ (>10 microns) (0.0000000882819%)

Carcinogen.

CAS: 14808-60-7

Density: 0

Relative Density: 2.6

Molecular Weight: 60.09

Boiling Point: 2230

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

POLYOXYETHYLENE (20) STEARYL ETHER (0.00000005331186%)

CAS: 9005-00-9

Density: 1

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 3 Quotient = 0.000

QUARTZ (<10 microns) (0.00000004414095%)

Carcinogen.

CAS: 14808-60-7

Density: 0

Relative Density: 2.6

Molecular Weight: 60.09  
Boiling Point: 2230  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 6 Quotient = 0.000  
FAD 3 Quotient = 0.000

ETHYLENE GLYCOL (0.0000000409801%)

Organic Solvent.  
CAS: 107-21-1  
Density: 1.11  
Relative Density: 1.1  
Molecular Weight: 62.07  
Boiling Point: 197.4  
Vapour Pressure: 0.05  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 2 Quotient = 0.000

ETHANOLAMINE (0.00000001927399%)

Organic Solvent.  
CAS: 141-43-5  
Density: 1.018  
Relative Density: 1.02  
Molecular Weight: 61.08  
Boiling Point: 170.8  
Vapour Pressure: 0.4  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 500. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 3 Quotient = 0.000  
FAD 2 Quotient = 0.000

SILANE,DICHLORODIMETHYL-,REACTION PRODUCTS WITH SILICA (0.00000001927399%)

CAS: 68611-44-9  
Density: 2  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 1 Quotient = 0.000

ALUMINUM OXIDE (0.00000000358833%)

CAS: 1344-28-1  
Density: 3.97  
Relative Density: 4  
Molecular Weight: 101.96  
Boiling Point: 3000  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 0. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.



FAD 1 Quotient = 0.000  
HYDROCHLORIC ACID (0.00000000239222%)  
CAS: 7647-01-0  
Density: 0.86  
Boiling Point: 109.85  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 2900. Limit: 0  
FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.  
FAD 4 Quotient = 0.000  
FAD 3 Quotient = 0.000

CYCLOHEXANE (0.0000000019274%)  
Organic Solvent.  
CAS: 110-82-7  
Density: 0.77  
Relative Density: 0.8  
Molecular Weight: 84.16  
Boiling Point: 80.7  
Vapour Pressure: 93.00791  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor entered: 13. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 0.000

ETHYL ALCOHOL (0.0000000019274%)  
Organic Solvent.  
CAS: 64-17-5  
Density: 0.786  
Relative Density: 0.8  
Molecular Weight: 46.08  
Boiling Point: 78.29  
Vapour Pressure: 42.94865  
LBLFactor = 200 (CAS=64175)  
MAL Factor entered: 7. Limit: 0  
FAD entered: 1; Lower Limit: 0  
FAD 1 Quotient = 0.000

2-PYRIDINETHIOL-1-OXIDE SODIUM SALT (0.000000006853%)  
CAS: 3811-73-2  
Density: 0  
Molecular Weight: 150.16  
Vapour Pressure: 0.00000034  
No LBL Factor entered or estimated from CAS Number or Boiling Point.  
MAL Factor from OEL: 0  
R Phrases: Xn;R22 Xn;R21 Xn;R20 Xi;R38 Xi;R36 N;R50  
FAD: 1. (Default)  
FAD 1 Quotient = 0.000

SODIUM NITRATE (0.000000005796%)  
CAS: 7631-99-4  
Density: 2.3  
Molecular Weight: 84.99

Boiling Point: 380

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

Diiron trioxide (0.00000000035883%)

CAS: 1309-37-1

Density: 5.25

Relative Density: 5.18

Molecular Weight: 159.7

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

SODIUM CHLORIDE (0.0000000002196%)

CAS: 7647-14-5

Density: 2.165

Molecular Weight: 58.44

Boiling Point: 1430.85

No LBL Factor entered or estimated from CAS Number or Boiling Point.

MAL Factor entered: 0. Limit: 0

FAD entered: 1; Lower Limit: No limit specified. A very low value will be used.

FAD 1 Quotient = 0.000

Density = 1.214. Entered value.

Figure-before-the dash = 00

WATER(@46.28%). MAL Factor = 0. Total increased by  $46.28*0=0$ . Running Total = 0

TITANIUM DIOXIDE(@18.03%). MAL Factor = 0. Total increased by  $18.03*0=0$ . Running Total = 0

2-(2-BUTOXYETHOXY)ETHANOL(@1.86%). MAL Factor = 0. Total increased by  $1.86*0=0$ . Running Total = 0

SODIUM POTASSIUM ALUMINUM SILICATE(@1.5%). MAL Factor = 0. Total increased by  $1.5*0=0$ . Running Total = 0

POLYPROPYLENE GLYCOL(@1.45%). MAL Factor = 0. Total increased by  $1.45*0=0.00$ . Running Total = 0.00

ALUMINUM HYDROXIDE(@0.76%). MAL Factor = 0. Total increased by  $0.76*0=0$ . Running Total = 0.00

POLYURETHANE RESIN(@0.39%). MAL Factor = 0. Total increased by  $0.39*0=0$ . Running Total = 0.00

Alcohols, C16-18 and C18-unsatd., ethoxylated(@0.34%). MAL Factor = 0. Total increased by  $0.34*0=0$ . Running Total = 0.00

Tripropylene glycol monomethyl ether(@0.29%). MAL Factor = 0. Total increased by  $0.29*0=0$ . Running Total = 0.00

Methanone, (1-hydroxycyclohexyl)phenyl-(@0.2%). MAL Factor = 0. Total increased by  $0.2*0=0$ . Running Total = 0.00

BENZOPHENONE(@0.20%). MAL Factor = 0. Total increased by  $0.20*0=0$ . Running Total = 0.00

DIPROPYLENE GLYCOL MONOMETHYL ETHER(@0.12%). MAL Factor = 5. Total increased by  $0.12*5=0.58$ . Running Total = 0.58

TRIMETHYLOLPROPANE(@0.09%). MAL Factor = 0. Total increased by  $0.09*0=0$ . Running Total = 0.58

SODIUM NITRITE(@0.06%). MAL Factor = 0. Total increased by  $0.06*0=0$ . Running Total = 0.58

3-Iodo-2-propynyl butylcarbamate(@0.06%). MAL Factor = 0. Total increased by  $0.06*0=0$ . Running Total = 0.58

ZIRCONIUM TETRAHYDROXIDE(@0.06%). MAL Factor = 0. Total increased by  $0.06*0=0$ . Running Total = 0.58

AMMONIUM HYDROXIDE(@0.05%). MAL Factor = 50. Total increased by  $0.05*50=2.29$ . Running Total = 2.87

AMMONIUM BENZOATE(@0.04%). MAL Factor = 0. Total increased by  $0.04*0=0$ . Running Total = 2.87

POLYETHYLENE-POLYPROPYLENE POLYMER(@0.04%). MAL Factor = 0. Total increased by  $0.04*0=0$ . Running Total = 2.87

1,2-BENZISOTHAZOLONE(@0.03%). MAL Factor = 0. Total increased by  $0.03*0=0$ . Running Total = 2.87

DIMETHYLAMINOETHANOL(@0.03%). MAL Factor = 280. Total increased by  $0.03*280=7.84$ . Running Total = 10.71

TITANIUM DIOXIDE (<10 microns)(@0.02%). MAL Factor = 0. Total increased by  $0.02*0=0$ . Running Total = 10.71

ZINC OXIDE(@0.02%). MAL Factor = 0. Total increased by  $0.02*0=0$ . Running Total = 10.71

2-BUTOXY ETHANOL(@0.01%). MAL Factor = 25. Total increased by  $0.01*25=0.14$ . Running Total = 10.85  
pyrithione zinc(@0.01%). MAL Factor = 0. Total increased by  $0.01*0=0$ . Running Total = 10.85  
1-BUTANOL(@0.00%). MAL Factor = 67. Total increased by  $0.00*67=0.12$ . Running Total = 10.97  
HEPTANE(@0.00%). MAL Factor = 12. Total increased by  $0.00*12=0.00$ . Running Total = 10.98  
DIETHYLENE GLYCOL(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ACETIC ACID(@0.00%). MAL Factor = 1. Total increased by  $0.00*1=0.00$ . Running Total = 10.98  
Triethyleneglycol monobutylether(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
PROPYLENE OXIDE(@0.00%). MAL Factor = 1. Total increased by  $0.00*1=0.00$ . Running Total = 10.98  
TRIETHYLENEGLYCOL(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
2-METHYL-4-ISOTHIAZOLIN-3-ONE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
OCTAMETHYLCYCLOTETRASILOXANE(@0.00%). MAL Factor = 1. Total increased by  $0.00*1=0.00$ . Running Total = 10.98  
Decamethylcyclopentasiloxane(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
alkyl polyglycol ether phosphate compound(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ETHANOL,2-(2-ETHOXYETHOXY)(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
Triethylene glycol monoethyl ether(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0.00$ . Running Total = 10.98  
ALUMINUM SILICATE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
AMORPHOUS SILICA(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
2-ETHYLHEXANOIC ACID(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
SILICA(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
LECITHINS(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
Ethanol, 2,2'-(butylimino)bis-(@0.00%). MAL Factor = 1. Total increased by  $0.00*1=0.00$ . Running Total = 10.98  
BENZENE(@0.00%). MAL Factor = 880. Total increased by  $0.00*880=0.00$ . Running Total = 10.98  
ZIRCONIUM OXIDE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
TIN(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0.00$ . Running Total = 10.98  
SODIUM HYDROXIDE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ACETALDEHYDE(@0.00%). MAL Factor = 1. Total increased by  $0.00*1=0.00$ . Running Total = 10.98  
2-BROMO-2-NITRO-1,3-PROPANEDIOL(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
1,4-DIOXANE(@0.00%). MAL Factor = 390. Total increased by  $0.00*390=0.00$ . Running Total = 10.98  
FORMALDEHYDE(@0.00%). MAL Factor = 2500. Total increased by  $0.00*2500=0.00$ . Running Total = 10.98  
METHYL ALCOHOL(@0.00%). MAL Factor = 54. Total increased by  $0.00*54=0.00$ . Running Total = 10.98  
ETHYLENE OXIDE(@0.00%). MAL Factor = 11. Total increased by  $0.00*11=0.00$ . Running Total = 10.98  
METHYL CHLORIDE(@0.00%). MAL Factor = 476.19. Total increased by  $0.00*476.19=0.00$ . Running Total = 10.98  
QUARTZ (>10 microns)(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
POLYOXYETHYLENE (20) STEARYL ETHER(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
QUARTZ (<10 microns)(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ETHYLENE GLYCOL(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ETHANOLAMINE(@0.00%). MAL Factor = 500. Total increased by  $0.00*500=0.00$ . Running Total = 10.98  
SILANE,DICHLORODIMETHYL-,REACTION PRODUCTS WITH SILICA(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
ALUMINUM OXIDE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
HYDROCHLORIC ACID(@0.00%). MAL Factor = 2900. Total increased by  $0.00*2900=0.00$ . Running Total = 10.98  
CYCLOHEXANE(@0.00%). MAL Factor = 13. Total increased by  $0.00*13=0.00$ . Running Total = 10.98  
ETHYL ALCOHOL(@0.00%). MAL Factor = 7. Total increased by  $0.00*7=0.00$ . Running Total = 10.98  
2-PYRIDINETHIOL-1-OXIDE SODIUM SALT(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0.00$ . Running Total = 10.98  
SODIUM NITRATE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
Diiron trioxide (@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
SODIUM CHLORIDE(@0.00%). MAL Factor = 0. Total increased by  $0.00*0=0$ . Running Total = 10.98  
Figure-before-the-dash calculated as 00. Via MAL Factor Total \* Density (10.98 \* 1.214) giving a MAL Number of 13  
MAL Number = Density (1.214) \* Sum (10.98) = 13

Figure-after-the-dash = 3. Calculated from component data.

proprietary acrylic copolymer (@26.32%) Increasing Total for FAD1 by 26322.82, giving 26322.82  
TITANIUM DIOXIDE (@18.03%) Increasing Total for FAD1 by 18034.3695773922, giving 44357.1895773922  
2-(2-BUTOXYETHOXY)ETHANOL (@1.86%) Increasing Total for FAD3 by 0.18632053, giving 0.18632053  
SODIUM POTASSIUM ALUMINUM SILICATE (@1.5%) Increasing Total for FAD1 by 15, giving 44372.1895773922  
POLYPROPYLENE GLYCOL (@1.45%) Increasing Total for FAD3 by 1452.5, giving 1452.68632053  
ALUMINUM HYDROXIDE (@0.76%) Increasing Total for FAD1 by 7.6002275, giving 44379.7898048922  
POLYMER, POLYFUNCTIONAL, NON-ANIONIC (@0.44%) Increasing Total for FAD1 by 441.9983, giving 44821.7881048922  
POLYURETHANE RESIN (@0.39%) Increasing Total for FAD1 by 3.94235, giving 44825.7304548922  
COALESCING AID (@0.35%) Increasing Total for FAD1 by 349.6479, giving 45175.3783548922  
Alcohols, C16-18 and C18-unsatd., ethoxylated (@0.34%) Increasing Total for FAD1 by 3.4023078279291, giving 45178.7806627201291  
POLYSILOXANE MIXTURE (@0.30%) Increasing Total for FAD1 by 297.362275, giving 45476.1429377201291  
Tripropylene glycol monomethyl ether (@0.29%) Increasing Total for FAD1 by 291.55, giving 45767.6929377201291  
proprietary surfactant (@0.28%) Increasing Total for FAD1 by 280.03, giving 46047.7229377201291  
POLYACRYLATE (@0.23%) Increasing Total for FAD1 by 229.8666, giving 46277.5895377201291  
Methanone, (1-hydroxycyclohexyl)phenyl- (@0.2%) Increasing Total for FAD1 by 2, giving 46279.5895377201291  
BENZOPHENONE (@0.20%) Increasing Total for FAD3 by 0.0998, giving 1452.78612053  
DIPROPYLENE GLYCOL MONOMETHYL ETHER (@0.12%) Increasing Total for FAD1 by 116.45, giving 46396.0395377201291  
TRIMETHYLOPROPANE (@0.09%) Increasing Total for FAD1 by 0.85312925, giving 46396.8926669701291  
SODIUM NITRITE (@0.06%) Increasing Total for FAD6 by 0.32, giving 0.32  
SODIUM NITRITE (@0.06%) Increasing Total for FAD3 by 0.64, giving 1453.42612053  
3-Iodo-2-propynyl butylcarbamate (@0.06%) Increasing Total for FAD3 by 0.0599421590592, giving 1453.4860626890592  
ZIRCONIUM TETRAHYDROXIDE (@0.06%) Increasing Total for FAD1 by 0.57, giving 46397.4626669701291  
reaction mass of mixed (3,3,4,4,5,5,6,6,7,7, 8,8,8- tridecafluorooctyl) phosphates, ammonium salt (@0.05%) Increasing Total for FAD1 by 54.4, giving 46451.8626669701291  
AMMONIUM HYDROXIDE (@0.05%) Increasing Total for FAD4 by 0.001306, giving 0.001306  
AMMONIUM HYDROXIDE (@0.05%) Increasing Total for FAD3 by 0.009142, giving 1453.4952046890592  
polyurethane resin (@0.04%) Increasing Total for FAD1 by 42.5, giving 46494.3626669701291  
AMMONIUM BENZOATE (@0.04%) Increasing Total for FAD3 by 0.04, giving 1453.5352046890592  
proprietary defoamer (@0.04%) Increasing Total for FAD1 by 39.2042, giving 46533.5668669701291  
POLYETHYLENE-POLYPROPYLENE POLYMER (@0.04%) Increasing Total for FAD1 by 0.3780342031032, giving 46533.9449011732323  
1,2-BENZISOTHIAZOLONE (@0.03%) Increasing Total for FAD3 by 0.0330321585497, giving 1453.5682368476089  
DIMETHYLAMINOETHANOL (@0.03%) Increasing Total for FAD3 by 0.0028, giving 1453.5710368476089  
DIMETHYLAMINOETHANOL (@0.03%) Increasing Total for FAD2 by 0.014, giving 0.014  
TITANIUM DIOXIDE (<10 microns) (@0.02%) Increasing Total for FAD1 by 16.245, giving 46550.1899011732323  
ZINC OXIDE (@0.02%) Increasing Total for FAD1 by 15, giving 46565.1899011732323  
polyethylene glycol monobutyl ether (@0.01%) Increasing Total for FAD1 by 9.25, giving 46574.4399011732323  
2-BUTOXY ETHANOL (@0.01%) Increasing Total for FAD3 by 0.00055823, giving 1453.5715950776089  
pyrithione zinc (@0.01%) Increasing Total for FAD3 by 0.0055, giving 1453.5770950776089  
1-BUTANOL (@0.00%) Increasing Total for FAD1 by 1.85, giving 46576.2899011732323  
ISOTHIAZOLONE SOLUTION (@0.00%) Increasing Total for FAD1 by 0.86624186471, giving 46577.1561430379423  
HEPTANE (@0.00%) Increasing Total for FAD1 by 0.396, giving 46577.5521430379423  
DIETHYLENE GLYCOL (@0.00%) Increasing Total for FAD3 by 0.000019057, giving 1453.5771141346089  
ACETIC ACID (@0.00%) Increasing Total for FAD4 by 0.00000752928196, giving 0.00131352928196  
Triethyleneglycol monobutylether (@0.00%) Increasing Total for FAD3 by 0.0000765, giving 1453.5771906346089  
Triethyleneglycol monobutylether (@0.00%) Increasing Total for FAD1 by 0.00153, giving 46577.5536730379423  
residual monomers (@0.00%) Increasing Total for FAD1 by 0.1334, giving 46577.6870730379423  
PROPYLENE OXIDE (@0.00%) Increasing Total for FAD6 by 0.0002625, giving 0.3202625

TRIETHYLENEGLYCOL (@0.00%) Increasing Total for FAD1 by 0.0004089609079, giving 46577.6874819988502  
2-METHYL-4-ISOTHIAZOLIN-3-ONE (@0.00%) Increasing Total for FAD6 by 0.00003588009408, giving 0.32029838009408  
2-METHYL-4-ISOTHIAZOLIN-3-ONE (@0.00%) Increasing Total for FAD3 by 0.001196003136, giving 1453.5783866377449  
OCTAMETHYLCYCLOTETRASILOXANE (@0.00%) Increasing Total for FAD3 by 0.0000280019274, giving 1453.5784146396723  
Decamethylcyclopentasiloxane (@0.00%) Increasing Total for FAD1 by 0.000280019274, giving 46577.6877620181242  
dodecamethylcyclohexasiloxane (@0.00%) Increasing Total for FAD1 by 0.0280019274, giving 46577.7157639455242  
alkyl polyglycol ether phosphate compound (@0.00%) Increasing Total for FAD3 by 0.000012337242405, giving 1453.578426976914705  
ETHANOL;2-(2-ETHOXYETHOXY) (@0.00%) Increasing Total for FAD3 by 0.00000153, giving 1453.578428506914705  
Triethylene glycol monoethyl ether (@0.00%) Increasing Total for FAD1 by 0.0153, giving 46577.7310639455242  
ALUMINUM SILICATE (@0.00%) Increasing Total for FAD1 by 0.0001458122715, giving 46577.7312097577957  
AMORPHOUS SILICA (@0.00%) Increasing Total for FAD1 by 0.0001195212918, giving 46577.7313292790875  
2-ETHYLHEXANOIC ACID (@0.00%) Increasing Total for FAD3 by 0.00000735, giving 1453.578435856914705  
SILICA (@0.00%) Increasing Total for FAD1 by 0.00650023922, giving 46577.7378295183075  
LECITHINS (@0.00%) Increasing Total for FAD1 by 0.0000528808926, giving 46577.7378823992001  
Ethanol, 2,2'-(butylimino)bis- (@0.00%) Increasing Total for FAD3 by 0.000002224425, giving 1453.578438081339705  
BENZENE (@0.00%) Increasing Total for FAD6 by 0.00004, giving 0.32033838009408  
ZIRCONIUM OXIDE (@0.00%) Increasing Total for FAD1 by 0.0000325, giving 46577.7379148992001  
TIN (@0.00%) Increasing Total for FAD1 by 0.0017, giving 46577.7396148992001  
POLYETHER SILOXANE COPOLYMER (@0.00%) Increasing Total for FAD1 by 0.00042488445, giving 46577.7400397836501  
SODIUM HYDROXIDE (@0.00%) Increasing Total for FAD4 by 0.0000004019274, giving 0.00131393120936  
SODIUM HYDROXIDE (@0.00%) Increasing Total for FAD3 by 0.000010048185, giving 1453.578448129524705  
ACETALDEHYDE (@0.00%) Increasing Total for FAD3 by 0.0000035, giving 1453.578451629524705  
polycarbonic acid ammonium salt (@0.00%) Increasing Total for FAD1 by 0.00030709786, giving 46577.7403468815101  
2-BROMO-2-NITRO-1,3-PROPANEDIOL (@0.00%) Increasing Total for FAD3 by 0.0000001988, giving 1453.578451828324705  
1,4-DIOXANE (@0.00%) Increasing Total for FAD6 by 0.0000000175, giving 0.32033839759408  
1,4-DIOXANE (@0.00%) Increasing Total for FAD3 by 0.00000175, giving 1453.578453578324705  
FORMALDEHYDE (@0.00%) Increasing Total for FAD6 by 0.000000175, giving 0.32033857259408  
FORMALDEHYDE (@0.00%) Increasing Total for FAD3 by 0.00000175, giving 1453.578455328324705  
METHYL ALCOHOL (@0.00%) Increasing Total for FAD6 by 0.00000000875, giving 0.32033858134408  
METHYL ALCOHOL (@0.00%) Increasing Total for FAD3 by 0.000000175, giving 1453.578455503324705  
ETHYLENE OXIDE (@0.00%) Increasing Total for FAD6 by 0.000000875, giving 0.32033945634408  
METHYL CHLORIDE (@0.00%) Increasing Total for FAD1 by 0.000175, giving 46577.7405218815101  
QUARTZ (>10 microns) (@0.00%) Increasing Total for FAD1 by 0.000000882819, giving 46577.7405227643291  
POLYOXYETHYLENE (20) STEARYL ETHER (@0.00%) Increasing Total for FAD3 by 0.00000002665593, giving 1453.578455529980635  
QUARTZ (<10 microns) (@0.00%) Increasing Total for FAD6 by 0.000000004414095, giving 0.320339460758175  
QUARTZ (<10 microns) (@0.00%) Increasing Total for FAD3 by 0.00000004414095, giving 1453.578455574121585  
ETHYLENE GLYCOL (@0.00%) Increasing Total for FAD2 by 0.00000000409801, giving 0.01400000409801  
ETHANOLAMINE (@0.00%) Increasing Total for FAD3 by 0.000000001927399, giving 1453.578455576048984  
ETHANOLAMINE (@0.00%) Increasing Total for FAD2 by 0.000000009636995, giving 0.014000013735005  
SILANE,DICHLORODIMETHYL-,REACTION PRODUCTS WITH SILICA (@0.00%) Increasing Total for FAD1 by 0.0000001927399, giving 46577.7405229570690  
ALUMINUM OXIDE (@0.00%) Increasing Total for FAD1 by 0.0000000358833, giving 46577.7405229929523  
HYDROCHLORIC ACID (@0.00%) Increasing Total for FAD4 by 0.00000000478444, giving 0.001313931687804  
HYDROCHLORIC ACID (@0.00%) Increasing Total for FAD3 by 0.00000000598055, giving 1453.578455582029534  
CYCLOHEXANE (@0.00%) Increasing Total for FAD1 by 0.0000019274, giving 46577.7405249203523  
ETHYL ALCOHOL (@0.00%) Increasing Total for FAD1 by 0.0000019274, giving 46577.7405268477523  
2-PYRIDINETHIOL-1-OXIDE SODIUM SALT (@0.00%) Increasing Total for FAD1 by 0.0000006853, giving 46577.7405275330523  
SODIUM NITRATE (@0.00%) Increasing Total for FAD1 by 0.000000005796, giving 46577.7405275388483  
Diiron trioxide (@0.00%) Increasing Total for FAD1 by 0.000000035883, giving 46577.7405275424366

SODIUM CHLORIDE (@0.00%) Increasing Total for FAD1 by 0.000000002196, giving 46577.7405275446326  
Figure-after-the-dash =3. Total of components with FAD=3 is >=1.

Low Boiling Liquid = False.

AMMONIUM HYDROXIDE (@0.05%) Total increased by  $0.05 \times 50 / 100 = 0.02$ . Running Total = 0.02

PROPYLENE OXIDE (@0.00%) Total increased by  $0.00 \times 1 / 100 = 0.00$ . Running Total = 0.02

ACETALDEHYDE (@0.00%) Total increased by  $0.00 \times 1 / 100 = 0.00$ . Running Total = 0.02

METHYL ALCOHOL (@0.00%) Total increased by  $0.00 \times 54 / 100 = 0.00$ . Running Total = 0.02

ETHYLENE OXIDE (@0.00%) Total increased by  $0.00 \times 11 / 100 = 0.00$ . Running Total = 0.02

METHYL CHLORIDE (@0.00%) Total increased by  $0.00 \times 476.19 / 100 = 0.00$ . Running Total = 0.02

ETHYL ALCOHOL (@0.00%) Total increased by  $0.00 \times 7 / 200 = 0.00$ . Running Total = 0.02

Density \* (Sum of components Concentration \* MALFactor/LBLFactor) = 0.03

Recommended Usage Temperature is < 40C, hence no MAL Code in use is assigned.

### Audit - RFU MAL Code

EU Denmark RFU MAL Code:-

Nothing was found

### New Fields for IA3.3

**MAL-code** : 00-3

**MAL Number** : 3.3302

**MAL Number (RFU)** : Not applicable.

**Protection based on MAL** : **According to the regulations on work involving coded products, the following stipulations apply to the use of personal protective equipment:**

**General:** Gloves must be worn for all work that may result in soiling. Apron/coveralls/protective clothing must be worn when soiling is so great that regular work clothes do not adequately protect skin against contact with the product. A face shield must be worn in work involving spattering if a full mask is not required. In this case, other recommended use of eye protection is not required.

In all spraying operations in which there is return spray, the following must be worn: respiratory protection and arm protectors/apron/coveralls/protective clothing as appropriate or as instructed.

MAL-code: 00-3

**Application:** During downtimes, cleaning and repair of closed facilities, spray booths or cabins, if there is a risk of contact with wet paint or organic solvents. When using scraper or knife, brush, roller, etc. for pre- and post-treatments in cabins or booths of the existing\* facility type, if the operator is inside the spray zone.

- Coveralls must be worn.

When spraying in existing\* spray booths, if the operator is outside the spray zone.

- Arm protectors and apron must be worn.

During all spraying where atomization occurs in cabins or spray booths where the

operator is inside the spray zone and during spraying outside a closed facility, cabin or booth.

- Air-supplied full mask, coveralls and hood must be worn.

**Drying:** Items for drying/drying ovens that are temporarily placed on such things as rack trolleys, etc. must be equipped with a mechanical exhaust system to prevent fumes from wet items from passing through workers' inhalation zone.

**Polishing:** When polishing treated surfaces, a mask with dust filter must be worn. When machine grinding, eye protection must be worn. Work gloves must always be worn.

**Caution** The regulations contain other stipulations in addition to the above.

\*See Regulations.

**Protection based on R-F-U MAL** : Not available.

Not available.

Not available.