# Audit - EU DK MAL Code

### **PPG VIKOTE 56 GREEN 4199**

## **Denmark MAL Code**

#### Audit - MAL Code

U Denmark MAL Code:- 5-3 The MAL Code calculations are performed with product and component data. Product is a Liquid PPG VIKOTE 56 GREEN 4199 - Components considered for the MAL Code calculation. {Denmark MAL Code} Hydrocarbons, C9, aromatics (31.98%) CAS: 64742-95-6 Density: 0.879 Molecular Weight: 123 Boiling Point: 172.5 Vapour Pressure: 1.5 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 58. Limit: 0 FAD entered: 1: Lower Limit: 0.1 FAD 1 Quotient = 319.8 ACRYLIC RESIN (28%) CAS: 25987-66-0 Density: 1.05 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 = 280 XYLENES (15.99225%) Organic Solvent. CAS: 1330-20-7 Density: 0.86 Relative Density: 0.861 Molecular Weight: 106.17 Boiling Point: 136.16 Vapour Pressure: 6.7 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 46. Limit: 0 FAD entered: 3; Lower Limit: 10 FAD 3 Quotient = 1.599FAD 1 Quotient = 79.961 SOLVENT NAPHTHA (PETROLEUM), LIGHT AROMATIC (8.4%) CAS: 64742-95-6 Density: 0.878 Molecular Weight: 123 Boiling Point: 172.5

Vapour Pressure: 1.500123 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 58. Limit: 0 FAD entered: 1: Lower Limit: 0.1 FAD 1 Quotient = 84 PARAFFIN WAXES AND HYDROCARBON WAXES; CHLORINATED (4.1%) CAS: 63449-39-8 Density: 1.21 Relative Density: 1 Molecular Weight: 462 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 = 4100 IRON HYDROXIDE OXIDE (3.997364%) CAS: 51274-00-1 Density: 4.26 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 = 39.974 ETHYLBENZENE (3.6530785%) Organic Solvent. Carcinogen. CAS: 100-41-4 Density: 0.866 Relative Density: 0.9 Molecular Weight: 106.18 Boiling Point: 136.1 Vapour Pressure: 9.30076 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 46. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 3 Quotient = 0.365COPPER PHTHALOCYANINE GREEN (1.39%) CAS: 1328-53-6 Density: 2.2 Vapour Pressure: 0.000009 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 = 1390 QUATERN.AM.CPS,BIS(HYDROGEN.TALLOW ALKYL)DIMET.-,BENTONITE (0.921595%) CAS: 68953-58-2 Density: 1.7 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 = 9.216N,N-1,6-HEXANEDIYLBIS (12-HYDROXY-OCTADECANEIMIDE) (0.4%) CAS: 55349-01-4 Density: 1.06 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 = 4ETHYL ALCOHOL (0.332486875%) 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 = 332.487 1-METHOXY-2-PROPYL ACETATE (0.207452578756%) Organic Solvent. CAS: 108-65-6 Density: 0.962 Relative Density: 0.96 Molecular Weight: 132.18 Boiling Point: 145.8 Vapour Pressure: 2.7 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 19. Limit: 0 FAD entered: 1; Lower Limit: 0 FAD 1 Quotient = 207.453 COPPER PHTALOCYANINE (0.1302892355%) CAS: 147-14-8 Density: 1.62 Molecular Weight: 576.1 Vapour Pressure: 0.000072 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 = 0.043 acrylic copolymer (0.11007584%) CAS: SUB110897 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 = 110.076

cyclohexanone (0.0993%) Organic Solvent. CAS: 108-94-1 Density: 0.946 Relative Density: 0.95 Molecular Weight: 98.14 Boiling Point: 154.3 Vapour Pressure: 3.75 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 70. Limit: 0 FAD entered: 1: Lower Limit: 0 FAD 1 Quotient = 99.3 CARBON BLACK (0.0696%) CAS: 1333-86-4 Density: 1.8 Relative Density: 1.95 Molecular Weight: 12.01 Boiling Point: 4200 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 6 Quotient = 0.003 FAD 3 Quotient = 0.007BLOCKED COPOLYMER (0.0675%) CAS: SUB100054 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 1 Quotient = 0.675 BLOCK COPOLYMER (0.03199144%) CAS: SUB101356 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 = 31.991 N-BUTYL ACETATE (0.02661308%) Organic Solvent. CAS: 123-86-4 Density: 0.881 Relative Density: 0.88 Molecular Weight: 116.18 Boiling Point: 126 Vapour Pressure: 11.25096 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 14. Limit: 0

FAD entered: 1: Lower Limit: 0 FAD 1 Quotient = 26.613QUARTZ (>10 microns) (0.019%) 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 = 0.19METHYL ALCOHOL (0.01752751896%) 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 = 0.001FAD 3 Quotient = 0.018 QUARTZ (<10 microns) (0.0094087125%) 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 = 0.001FAD 3 Quotient = 0.009 DIMETHYL GLUTARATE (0.00920509212%) CAS: 1119-40-0 Density: 1.09 Molecular Weight: 160.17 Boiling Point: 216 Vapour Pressure: 0.062 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 4. Limit: 0 FAD entered: 1; Lower Limit: 0 FAD 1 Quotient = 9.205

copper phthalocyanine derivative (0.0066825%) CAS: SUB142534 Density: 0 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 = 6.682Copper, [29H,31H-phthalocyaninato(2-)-N29,N30,N31,N32]-, (1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)methyl derivs. (0.00594%) CAS: 68411-06-3 Density: 1.6 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 2 Quotient = 0.002N,N"-naphthalene-1,5-diylbis[N'-[3-[(2-ethylhexyl)oxy]propyl]urea] (0.004%) CAS: 71216-01-8 Density: 0 Molecular Weight: 584.83 No LBL Factor entered or estimated from CAS Number or Boiling Point. No MAL Factor calculated. FAD: 1. (Default) FAD 1 Quotient = 4 [[2,2',2"-[29H,31H-phthalocyaninetrivltris(methylene)]tris[1H-isoindole-1,3(2H)-dionato]](2-)-N29,N30,N31,N32]copper (0.004%) CAS: 59160-79-1 Density: 0 Molecular Weight: 1053.49 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 = 0.001 DIMETHYL SUCCINATE (0.00315092092%) CAS: 106-65-0 Density: 1.119 Molecular Weight: 146.16 Boiling Point: 196.2 Vapour Pressure: 0.18 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 = 3.151[N,N,N',N'',N'',N''-hexaethyl-29H,31H-phthalocyaninetrimethylaminato(2-)-N29,N30,N31,N32]copper tris(dodecylbenzenesulphonate) (0.00232%) CAS: 75247-18-6 Density: 0 Molecular Weight: 1810.99 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 = 0.001DIMETHYL ADIPATE (0.00136906826%) CAS: 627-93-0 Density: 1.062 Molecular Weight: 174.22 Boiling Point: 230.9 Vapour Pressure: 0.021 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 = 0.014METHYL METHACRYLATE (0.000982623436%) Organic Solvent. CAS: 80-62-6 Density: 0.94 Relative Density: 0.94 Molecular Weight: 100.13 Boiling Point: 100.36 Vapour Pressure: 27.75236 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 46. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 5 Quotient = 0.000 FAD 3 Quotient = 0.0012-METHOXY-1-PROPYL ACETATE (0.000841146496%) Organic Solvent. CAS: 70657-70-4 Density: 0.97 Molecular Weight: 132.18 Boiling Point: 150.5 Vapour Pressure: 2.9 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 181. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 6 Quotient = 0.004 ALUMINUM SILICATE (0.0007387875%) 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 = 0.007Siloxanes and Silicones, methyl 3,3,3-trifluoropropyl (0.0006999%) CAS: 63148-56-1 Density: 0 No LBL Factor entered or estimated from CAS Number or Boiling Point. No MAL Factor calculated.

FAD: 1. (Default) FAD 1 Quotient = 0.700MANGANESE (0.00068%) CAS: 7439-96-5 Density: 7.47 Molecular Weight: 54.94 Boiling Point: 1962 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 2 Quotient = 0.001 SILVER (0.000504%) CAS: 7440-22-4 Density: 10.49 Relative Density: 10.5 Molecular Weight: 107.87 Boiling Point: 2212 Vapour Pressure: 0.001 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 = 0.0052-Propenoic acid, 2-methyl-, 1,7,7-trimethylbicyclo[2.2.1]hept-2-yl ester, exo- (0.000471557036%) CAS: 7534-94-3 Density: 0.983 Molecular Weight: 222.33 Boiling Point: 275 Vapour Pressure: 0.009 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 5 Quotient = 0.000FAD 3 Quotient = 0.000N-BUTYL METHACRYLATE (0.000467036064%) Organic Solvent. CAS: 97-88-1 Density: 0.89 Relative Density: 0.9 Molecular Weight: 142.22 Boiling Point: 163 Vapour Pressure: 1.59014 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 16. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 5 Quotient = 0.000CHROMIUM (0.000436%) CAS: 7440-47-3

Density: 7.15 Relative Density: 7.14 Molecular Weight: 52 Boiling Point: 2642 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 = 0.000COPPER (0.000388%) CAS: 7440-50-8 Density: 8.78 Relative Density: 8.9 Molecular Weight: 63.55 Boiling Point: 2595 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 dodecyltrimethylammonium chloride (0.000274725%) CAS: 112-00-5 Density: 0 Molecular Weight: 263.95 No LBL Factor entered or estimated from CAS Number or Boiling Point. No MAL Factor calculated. FAD: 1. (Default) FAD 1 Quotient = 0.275 PROPYLENE GLYCOL MONOMETHYL ETHER (0.00023796296%) Organic Solvent. CAS: 107-98-2 Density: 0.92 Relative Density: 0.92 Molecular Weight: 90.14 Boiling Point: 120.17 Vapour Pressure: 8.5 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 28. Limit: 0 FAD entered: 1; Lower Limit: 0 FAD 1 Quotient = 0.238ZINC (0.000236%) CAS: 7440-66-6 Density: 7.1 Relative Density: 7.14 Molecular Weight: 65.37 Boiling Point: 908 Vapour Pressure: 0.00000075 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 = 0.002NICKEL (0.000184%) Carcinogen. CAS: 7440-02-0 Density: 8.9 Relative Density: 8.9 Molecular Weight: 58.71 Boiling Point: 2730 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 5 Quotient = 0.002FAD 6 Quotient = 0.000 organotin compound (0.0001485%) CAS: SUB143296 Density: 0 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 = 0.148TOLUENE (0.00009%) Organic Solvent. CAS: 108-88-3 Density: 0.87 Relative Density: 0.87 Molecular Weight: 92.14 Boiling Point: 110.6 Vapour Pressure: 23.17 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor entered: 74. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 3 Quotient = 0.000 MOLYBDENUM (0.00008%) CAS: 7439-98-7 Density: 10.2 Relative Density: 10.28 Molecular Weight: 95.94 Boiling Point: 4612 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 = 0.08 WATER (0.00007457878%) 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 2-TERT-BUTYLAMINOETHYL METHACRYLATE (0.0000393128%) CAS: 3775-90-4 Density: 0.914 Relative Density: 0.9 Molecular Weight: 185.3 Boiling Point: 215 Vapour Pressure: 0.04 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 = 0.000FAD 5 Quotient = 0.000 ANTIMONY (0.000036%) CAS: 7440-36-0 Density: 6.7 Molecular Weight: 121.75 Boiling Point: 1635 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor from OEL: 0 R Phrases: T:R25 FAD: 1. (Default) FAD 1 Quotient = 0.036 CUMENE (0.00003%) Organic Solvent. CAS: 98-82-8 Density: 0.86 Relative Density: 0.9 Molecular Weight: 120.21 Boiling Point: 152 Vapour Pressure: 3.72032 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 = 0.000BARIUM (0.000028%) CAS: 7440-39-3 Density: 3.6 Relative Density: 3.6 Molecular Weight: 137.34 Boiling Point: 1640 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor from OEL: 0 R Phrases: F;R15 Xi;R38 Xi;R36 Xi;R37

FAD: 1. (Default) FAD 1 Quotient = 0.0281-BUTANOL (0.00002583924%) 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 = 0.026ARSENIC (0.000024%) Carcinogen. CAS: 7440-38-2 Density: 5.7 Relative Density: 5.73 Molecular Weight: 74.92 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 6 Quotient = 0.000 ACETIC ACID (0.00001923353%) 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: 400. Limit: 0 FAD entered: 1; Lower Limit: No limit specified. A very low value will be used. FAD 4 Quotient = 0.000FAD 3 Quotient = 0.000 COBALT (0.000012%) Carcinogen. CAS: 7440-48-4 Density: 8.9 Relative Density: 8.92 Molecular Weight: 58.93 Boiling Point: 2870 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 = 0.000VANADIUM (0.000012%) CAS: 7440-62-2 Density: 6.57 Relative Density: 6.11 Molecular Weight: 50.94 Boiling Point: 3000 No LBL Factor entered or estimated from CAS Number or Boiling Point. No MAL Factor calculated. FAD: 1. (Default) FAD 1 Quotient = 0.012Lead (0.00008%) CAS: 7439-92-1 Density: 11.34 Relative Density: 11.34 Molecular Weight: 207.19 Boiling Point: 660 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 = 0.000FAD 6 Quotient = 0.000 CADMIUM (0.00008%) Carcinogen. CAS: 7440-43-9 Density: 8.64 Relative Density: 8.64 Molecular Weight: 112.4 Boiling Point: 766.85 Vapour Pressure: 0.97507995 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 = 0.000BUTYLATED HYDROXYTOLUENE (0.0000073548%) CAS: 128-37-0 Density: 1.03 Relative Density: 1.048 Molecular Weight: 220.39 Boiling Point: 265 Vapour Pressure: 0.00825 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 ISOBUTYL METHACRYLATE (0.000004717536%) Organic Solvent. CAS: 97-86-9

Density: 0.88 Relative Density: 0.8858 Molecular Weight: 142.22 Boiling Point: 155 Vapour Pressure: 1.58263 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 FAD 5 Quotient = 0.000BENZENE (0.0000045%) 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 DENATONIUM BENZOATE (0.000003325%) CAS: 3734-33-6 Density: 0 Molecular Weight: 446.59 No LBL Factor entered or estimated from CAS Number or Boiling Point. No MAL Factor calculated. FAD: 1. (Default) FAD 1 Quotient = 0.003 TIN (0.00000676742%) 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 = 0.001 POLYCHLOROBIPHENYLS (0.00000066825%) Carcinogen. CAS: 1336-36-3 Density: 0 No LBL Factor entered or estimated from CAS Number or Boiling Point. MAL Factor from OEL: 0

R Phrases: R33 N;R50/53 FAD: 1. (Default) FAD 1 Quotient = 0.001 ACETONE (0.000000525%) Organic Solvent. CAS: 67-64-1 Density: 0.791 Relative Density: 0.8 Molecular Weight: 58.09 Boiling Point: 56.05 Vapour Pressure: 180.01463 LBLFactor = 100 (BP=56.05) MAL Factor entered: 23. Limit: 0 FAD entered: 1: Lower Limit: 0 FAD 1 Quotient = 0.001N,N-Dimethyl-1-Aminododecane (0.00000037125%) CAS: 112-18-5 Density: 0.778 Molecular Weight: 213.46 Boiling Point: 260 Vapour Pressure: 0.00017 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 = 0.0004-METHOXYPHENOL (0.000000196564%) CAS: 150-76-5 Density: 1.6 Relative Density: 1.55 Molecular Weight: 124.15 Boiling Point: 243 Vapour Pressure: 0.00675 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 5 Quotient = 0.000OCTAMETHYLCYCLOTETRASILOXANE (0.0000001%) 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 Density = 1.005. Entered value.

Figure-before-the dash = 5Hydrocarbons, C9, aromatics (@31.98%). MAL Factor = 58. Total increased by 31.98\*58=1854.84. Running Total = 1854.84 ACRYLIC RESIN(@28%), MAL Factor = 0. Total increased by 28\*0=0. Running Total = 1854.84 XYLENES(@15.99%). MAL Factor = 46. Total increased by 15.99\*46=735.64. Running Total = 2590.48 SOLVENT NAPHTHA (PETROLEUM), LIGHT AROMATIC(@8.4%). MAL Factor = 58. Total increased by 8.4\*58=487.2. Running Total = 3077.68 PARAFFIN WAXES AND HYDROCARBON WAXES; CHLORINATED(@4.1%). MAL Factor = 0. Total increased by 4.1\*0=0. Running Total = 3077.68 IRON HYDROXIDE OXIDE(@4.00%). MAL Factor = 0. Total increased by 4.00\*0=0. Running Total = 3077.68 ETHYLBENZENE(@3.65%), MAL Factor = 46. Total increased by 3.65\*46=168.04. Running Total = 3245.73 COPPER PHTHALOCYANINE GREEN(@1.39%). MAL Factor = 0. Total increased by 1.39\*0=0. Running Total = 3245.73 QUATERN.AM.CPS,BIS(HYDROGEN.TALLOW ALKYL)DIMET.-,BENTONITE(@0.92%). MAL Factor = 0. Total increased by 0.92\*0=0. Running Total = 3245.73 N.N-1,6-HEXANEDIYLBIS (12-HYDROXY-OCTADECANEIMIDE)(@0.4%). MAL Factor = 0. Total increased by 0.4\*0=0. Running Total = 3245.73 ETHYL ALCOHOL(@0.33%). MAL Factor = 7. Total increased by 0.33\*7=2.33. Running Total = 3248.05 1-METHOXY-2-PROPYL ACETATE(@0.21%), MAL Factor = 19, Total increased by 0.21\*19=3.94, Running Total = 3251.99 COPPER PHTALOCYANINE(@0.13%). MAL Factor = 0. Total increased by 0.13\*0=0. Running Total = 3251.99 cyclohexanone(@0.10%). MAL Factor = 70. Total increased by 0.10\*70=6.95. Running Total = 3258.95 CARBON BLACK(@0.07%). MAL Factor = 0. Total increased by 0.07\*0=0. Running Total = 3258.95 BLOCKED COPOLYMER(@0.07%). MAL Factor = 0. Total increased by 0.07\*0=0. Running Total = 3258.95 N-BUTYL ACETATE(@0.03%), MAL Factor = 14. Total increased by 0.03\*14=0.37. Running Total = 3259.32 QUARTZ (>10 microns)(@0.02%). MAL Factor = 0. Total increased by 0.02\*0=0. Running Total = 3259.32 METHYL ALCOHOL(@0.02%). MAL Factor = 54. Total increased by 0.02\*54=0.95. Running Total = 3260.26 QUARTZ (<10 microns)(@0.01%). MAL Factor = 0. Total increased by 0.01\*0=0. Running Total = 3260.26 DIMETHYL GLUTARATE(@0.01%). MAL Factor = 4. Total increased by 0.01\*4=0.04. Running Total = 3260.30 copper phthalocyanine derivative(@0.01%). MAL Factor = 0. Total increased by 0.01\*0=0.00. Running Total = 3260.30 Copper, [29H,31H-phthalocyaninato(2-)-N29,N30,N31,N32]-, (1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)methyl derivs.(@0.01%). MAL Factor = 0. Total increased by 0.01\*0=0. Running Total = 3260.30 [[2,2',2"-[29H,31H-phthalocyaninetrivitris(methylene)]tris[1H-isoindole-1,3(2H)-dionato]](2-)-N29,N30,N31,N32]copper(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.30 DIMETHYL SUCCINATE(@0.00%). MAL Factor = 5. Total increased by 0.00\*5=0.02. Running Total = 3260.32 [N,N,N',N'',N'',N''-hexaethyl-29H,31H-phthalocyaninetrimethylaminato(2-)-N29,N30,N31,N32]copper tris(dodecylbenzenesulphonate)(@0.00%). MAL Factor = 0. Total increased by  $0.00^{\circ}0=0$ . Running Total = 3260.32 DIMETHYL ADIPATE(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.32 METHYL METHACRYLATE(@0.00%). MAL Factor = 46. Total increased by 0.00\*46=0.05. Running Total = 3260.36 2-METHOXY-1-PROPYL ACETATE(@0.00%). MAL Factor = 181. Total increased by 0.00\*181=0.15. Running Total = 3260.51 ALUMINUM SILICATE(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.51 MANGANESE(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.51 SILVER(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.51 2-Propenoic acid, 2-methyl-, 1,7,7-trimethylbicyclo[2.2.1]hept-2-yl ester, exo-(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.51 N-BUTYL METHACRYLATE(@0.00%). MAL Factor = 16. Total increased by 0.00\*16=0.01. Running Total = 3260.52 CHROMIUM(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.52 COPPER(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.52 PROPYLENE GLYCOL MONOMETHYL ETHER(@0.00%). MAL Factor = 28. Total increased by 0.00\*28=0.01. Running Total = 3260.53 ZINC(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.53 NICKEL(@0.00%). MAL Factor = 0. Total increased by  $0.00^{\circ}0=0$ . Running Total = 3260.53 organotin compound(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0.00. Running Total = 3260.53 TOLUENE(@0.00%). MAL Factor = 74. Total increased by 0.00\*74=0.01. Running Total = 3260.54 MOLYBDENUM(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0.00. Running Total = 3260.54 WATER(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 2-TERT-BUTYLAMINOETHYL METHACRYLATE(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54

ANTIMONY(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0.00. Running Total = 3260.54 CUMENE(@0.00%). MAL Factor = 1. Total increased by 0.00\*1=0.00. Running Total = 3260.54 BARIUM(@0.00%), MAL Factor = 0. Total increased by 0.00\*0=0.00, Running Total = 3260.54 1-BUTANOL(@0.00%), MAL Factor = 67, Total increased by 0.00\*67=0.00, Running Total = 3260.54 ARSENIC(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 ACETIC ACID(@0.00%). MAL Factor = 400. Total increased by 0.00\*400=0.01. Running Total = 3260.54 COBALT(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 Lead(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 CADMIUM(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 BUTYLATED HYDROXYTOLUENE(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.54 ISOBUTYL METHACRYLATE(@0.00%). MAL Factor = 1. Total increased by 0.00\*1=0.00. Running Total = 3260.54 BENZENE(@0.00%). MAL Factor = 880. Total increased by 0.00\*880=0.00. Running Total = 3260.55 TIN(@0.00%), MAL Factor = 0. Total increased by 0.00\*0=0.00. Running Total = 3260.55 POLYCHLOROBIPHENYLS(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0.00. Running Total = 3260.55 ACETONE(@0.00%). MAL Factor = 23. Total increased by 0.00\*23=0.00. Running Total = 3260.55 N,N-Dimethyl-1-Aminododecane(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.55 4-METHOXYPHENOL(@0.00%). MAL Factor = 0. Total increased by 0.00\*0=0. Running Total = 3260.55 OCTAMETHYLCYCLOTETRASILOXANE(@0.00%), MAL Factor = 1, Total increased by 0.00\*1=0.00, Running Total = 3260.55 Figure-before-the-dash calculated as 5. Via MAL Factor Total \* Density (3260.55 \* 1.005) giving a MAL Number of 3277 MAL Number = Density (1.005) \* Sum (3260.55) = 3277 Figure-after-the-dash = 3. Calculated from component data. Hydrocarbons, C9, aromatics (@31.98%) Increasing Total for FAD1 by 319.8, giving 319.8 ACRYLIC RESIN (@28%) Increasing Total for FAD1 by 280, giving 599.8 XYLENES (@15.99%) Increasing Total for FAD3 by 1.599225, giving 1.599225 XYLENES (@15.99%) Increasing Total for FAD1 by 79.96125, giving 679.76125 SOLVENT NAPHTHA (PETROLEUM), LIGHT AROMATIC (@8.4%) Increasing Total for FAD1 by 84, giving 763.76125 PARAFFIN WAXES AND HYDROCARBON WAXES; CHLORINATED (@4.1%) Increasing Total for FAD1 by 4100, giving 4863.76125 IRON HYDROXIDE OXIDE (@4.00%) Increasing Total for FAD1 by 39.97364, giving 4903.73489 ETHYLBENZENE (@3.65%) Increasing Total for FAD3 by 0.36530785, giving 1.96453285 COPPER PHTHALOCYANINE GREEN (@1.39%) Increasing Total for FAD1 by 1390, giving 6293,73489 QUATERN.AM.CPS,BIS(HYDROGEN.TALLOW ALKYL)DIMET.-,BENTONITE (@0.92%) Increasing Total for FAD1 by 9.21595, giving 6302.95084 N.N-1,6-HEXANEDIYLBIS (12-HYDROXY-OCTADECANEIMIDE) (@0.4%) Increasing Total for FAD1 by 4, giving 6306.95084 ETHYL ALCOHOL (@0.33%) Increasing Total for FAD1 by 332.486875, giving 6639.437715 1-METHOXY-2-PROPYL ACETATE (@0.21%) Increasing Total for FAD1 by 207.452578756, giving 6846.890293756 acrylic copolymer (@0.11%) Increasing Total for FAD1 by 110.07584, giving 6956.966133756 cyclohexanone (@0.10%) Increasing Total for FAD1 by 99.3, giving 7056.266133756 CARBON BLACK (@0.07%) Increasing Total for FAD6 by 0.002784, giving 0.002784 CARBON BLACK (@0.07%) Increasing Total for FAD3 by 0.00696, giving 1.97149285 BLOCKED COPOLYMER (@0.07%) Increasing Total for FAD1 by 0.675, giving 7056.941133756 BLOCK COPOLYMER (@0.03%) Increasing Total for FAD1 by 31.99144, giving 7088.932573756 N-BUTYL ACETATE (@0.03%) Increasing Total for FAD1 by 26.61308, giving 7115.545653756 QUARTZ (>10 microns) (@0.02%) Increasing Total for FAD1 by 0.19, giving 7115.735653756 METHYL ALCOHOL (@0.02%) Increasing Total for FAD6 by 0.000876375948, giving 0.003660375948 METHYL ALCOHOL (@0.02%) Increasing Total for FAD3 by 0.01752751896, giving 1.98902036896 QUARTZ (<10 microns) (@0.01%) Increasing Total for FAD6 by 0.00094087125, giving 0.004601247198 QUARTZ (<10 microns) (@0.01%) Increasing Total for FAD3 by 0.0094087125, giving 1.99842908146 DIMETHYL GLUTARATE (@0.01%) Increasing Total for FAD1 by 9.20509212, giving 7124.940745876

copper phthalocyanine derivative (@0.01%) Increasing Total for FAD1 by 6.6825, giving 7131.623245876

N,N"-naphthalene-1,5-diylbis[N'-[3-[(2-ethylhexyl)oxy]propyl]urea] (@0.00%) Increasing Total for FAD1 by 4, giving 7135.623245876

DIMETHYL SUCCINATE (@0.00%) Increasing Total for FAD1 by 3.15092092, giving 7138.774166796

DIMETHYL ADIPATE (@0.00%) Increasing Total for FAD1 by 0.0136906826, giving 7138.7878574786

METHYL METHACRYLATE (@0.000982623436%) Increasing Total for FAD5 by 0.0001965246872, giving 0.0001965246872

METHYL METHACRYLATE (@0.00%) Increasing Total for FAD3 by 0.000982623436, giving 1.999411704896

2-METHOXY-1-PROPYL ACETATE (@0.00%) Increasing Total for FAD6 by 0.00420573248, giving 0.008806979678

ALUMINUM SILICATE (@0.00%) Increasing Total for FAD1 by 0.007387875, giving 7138.7952453536

Siloxanes and Silicones, methyl 3,3,3-trifluoropropyl (@0.00%) Increasing Total for FAD1 by 0.6999, giving 7139.4951453536

SILVER (@0.00%) Increasing Total for FAD1 by 0.00504, giving 7139.5001853536

2-Propenoic acid, 2-methyl-, 1,7,7-trimethylbicyclo[2.2.1]hept-2-yl ester, exo- (@0.000471557036%) Increasing Total for FAD5 by 0.0000943114072, giving 0.0002908360944

2-Propenoic acid, 2-methyl-, 1,7,7-trimethylbicyclo[2.2.1]hept-2-yl ester, exo- (@0.00%) Increasing Total for FAD3 by 0.000471557036, giving 1.999883261932 N-BUTYL METHACRYLATE (@0.000467036064%) Increasing Total for FAD5 by 0.000467036064, giving 0.0007578721584

CHROMIUM (@0.00%) Increasing Total for FAD3 by 0.0000436, giving 1.999926861932

dodecyltrimethylammonium chloride (@0.00%) Increasing Total for FAD1 by 0.274725, giving 7139.7749103536

PROPYLENE GLYCOL MONOMETHYL ETHER (@0.00%) Increasing Total for FAD1 by 0.23796296, giving 7140.0128733136

ZINC (@0.00%) Increasing Total for FAD1 by 0.00236, giving 7140.0152333136

NICKEL (@0.00%) Increasing Total for FAD6 by 0.0000368, giving 0.008843779678

NICKEL (@0.000184%) Increasing Total for FAD5 by 0.00184, giving 0.0025978721584

organotin compound (@0.00%) Increasing Total for FAD1 by 0.1485, giving 7140.1637333136

TOLUENE (@0.00%) Increasing Total for FAD3 by 0.000009, giving 1.999935861932

MOLYBDENUM (@0.00%) Increasing Total for FAD1 by 0.08, giving 7140.2437333136

2-TERT-BUTYLAMINOETHYL METHACRYLATE (@0.0000393128%) Increasing Total for FAD5 by 0.00000786256, giving 0.0026057347184

2-TERT-BUTYLAMINOETHYL METHACRYLATE (@0.00%) Increasing Total for FAD3 by 0.0000393128, giving 1.999975174732

ANTIMONY (@0.00%) Increasing Total for FAD1 by 0.036, giving 7140.2797333136

CUMENE (@0.00%) Increasing Total for FAD3 by 0.00003, giving 2.000005174732

BARIUM (@0.00%) Increasing Total for FAD1 by 0.028, giving 7140.3077333136

1-BUTANOL (@0.00%) Increasing Total for FAD1 by 0.02583924, giving 7140.3335725536

ARSENIC (@0.00%) Increasing Total for FAD6 by 0.00012, giving 0.008963779678

ACETIC ACID (@0.00%) Increasing Total for FAD4 by 0.0000007693412, giving 0.0000007693412

ACETIC ACID (@0.00%) Increasing Total for FAD3 by 0.000001923353, giving 2.000007098085

COBALT (@0.00%) Increasing Total for FAD6 by 0.00012, giving 0.009083779678

VANADIUM (@0.00%) Increasing Total for FAD1 by 0.012, giving 7140.3455725536

Lead (@0.00%) Increasing Total for FAD6 by 0.0000008, giving 0.009084579678

Lead (@0.00%) Increasing Total for FAD3 by 0.000032, giving 2.000039098085

CADMIUM (@0.00%) Increasing Total for FAD6 by 0.00008, giving 0.009164579678

BUTYLATED HYDROXYTOLUENE (@0.00%) Increasing Total for FAD3 by 0.00000073548, giving 2.000039833565

ISOBUTYL METHACRYLATE (@0.000004717536%) Increasing Total for FAD5 by 0.0000009435072, giving 0.0026066782256

ISOBUTYL METHACRYLATE (@0.00%) Increasing Total for FAD3 by 0.000004717536, giving 2.000044551101

BENZENE (@0.00%) Increasing Total for FAD6 by 0.000045, giving 0.009209579678 DENATONIUM BENZOATE (@0.00%) Increasing Total for FAD1 by 0.003325, giving 7140.3488975536 TIN (@0.00%) Increasing Total for FAD1 by 0.000676742, giving 7140.3495742956 POLYCHLOROBIPHENYLS (@0.00%) Increasing Total for FAD1 by 0.00066825, giving 7140.3502425456 ACETONE (@0.00%) Increasing Total for FAD1 by 0.000525, giving 7140.3507675456 N,N-Dimethyl-1-Aminododecane (@0.00%) Increasing Total for FAD4 by 0.00000037125, giving 0.0000011405912 4-METHOXYPHENOL (@0.00000196564%) Increasing Total for FAD5 by 0.000000196564, giving 0.0026068747896 OCTAMETHYLCYCLOTETRASILOXANE (@0.00%) Increasing Total for FAD3 by 0.0000001, giving 2.000044651101 Figure-after-the-dash =3. Total of components with FAD=3 is >=1. Low Boiling Liquid = False.

ETHYL ALCOHOL (@0.33%) Total increased by 0.33\*7/200=0.01. Running Total = 0.01 METHYL ALCOHOL (@0.02%) Total increased by 0.02\*54/100=0.01. Running Total = 0.02 ACETONE (@0.00%) Total increased by 0.00\*23/100=0.00. Running Total = 0.02 Density \* (Sum of components Concentration \* MALFactor/LBLFactor) = 0.02 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	: 5-3
MAL Number	: 3276.85
MAL Number (RFU)	: Not applicable.
Protection based on MAL	: According to the

: 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: 5-3

**Application:** When spraying in new\* booths if the operator is outside the spray zone. During non-atomizing spraying in existing\* facilities of the combined-cabin, spray-cabin and spray-booth type where the operator is working inside the spray zone. When using scraper or knife, brush, roller, etc. for pre- and post-treatments outside a closed facility, spray booth or spray cabin.

- Air-supplied full mask must be worn.

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. 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.

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

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

- Air-supplied full mask, 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.

: Not available.

Protection based on R-F-U

MAL

Not available. Not available.