Declaration of Compliance

Nuvoton Technology Corporation hereby make declaration to our best knowledge that our products are free of SVHC (Substance of Very High Concern) of European Union REACH directive as listed below.

Item	Substance of Very High Concern	CAS number	EC number	Limit
1	Anthracene	120-12-7	204-371-1	< 0.1 %
2	4,4'- Diaminodiphenylmethane (MDA)	101-77-9	202-974-4	< 0.1 %
3	Dibutyl phthalate (DBP)	84-74-2	201-557-4	< 0.1 %
4	Cobalt dichloride	7646-79-9	231-589-4	< 0.1 %
5	Diarsenic pentaoxide	1303-28-2	215-116-9	< 0.1 %
6	Diarsenic trioxide	1327-53-3	215-481-4	< 0.1 %
7	Sodium dichromate	7789-12-0 10588-01-9	234-190-3	< 0.1 %
8	5-tert-butyl-2,4,6-trinitro-m-xylene (Musk xylene)	81-15-2	201-329-4	< 0.1 %
9	Bis (2-ethyl(hexyl)phthalate) (DEHP)	117-81-7	204-211-0	< 0.1 %
10	Hexabromocyclododecane (HBCDD) and all major diastereoisomers identified: Alpha-hexabromocyclododecane Beta-hexabromocyclododecane Gamma-hexabromocyclododecane	25637-99-4 3194-55-6 134237-50-6 134237-51-7 134237-52-8	247-148-4 221-695-9	< 0.1 %
11	Alkanes, C10-13, chloro (Short Chain Chlorinated Paraffins)	85535-84-8	287-476-5	< 0.1 %
12	Bis(tributyltin) oxide (TBTO)	56-35-9	200-268-0	< 0.1 %
13	Lead hydrogen arsenate	7784-40-9	232-064-2	< 0.1 %
14	Triethyl arsenate	15606-95-8	427-700-2	< 0.1 %
15	Benzyl butyl phthalate (BBP)	85-68-7	201-622-7	< 0.1 %
16	Anthracene oil	90640-80-5	292-602-7	< 0.1 %
17	Anthracene oil, anthracene paste, distn. Lights	91995-17-4	295-278-5	< 0.1 %
18	Anthracene oil, anthracene paste, anthracene fraction	91995-15-2	295-275-9	< 0.1 %
19	Anthracene oil, anthracene-low	90640-82-7	292-604-8	< 0.1 %
20	Anthracene oil, anthracene paste	90640-81-6	292-603-2	< 0.1 %
21	Diisobutyl phthalate (DIBP)	84-69-5	201-553-2	< 0.1 %
22	2,4-Dinitrotoluene	121-14-2	204-450-0	< 0.1 %
23	Coal tar pitch, high temperature	65996-93-2	266-028-2	< 0.1 %
24	Tris(2-chloroethyl)phosphate (TCEP)	115-96-8	204-118-5	< 0.1 %
25	Acrylamide	79-06-1	201-173-7	< 0.1 %
26	Lead chromate	7758-97-6	231-846-0	< 0.1 %

27	Lead chromate molybdate sulphate red (C.I. Pigment Red 104)	12656-85-8	235-759-9	< 0.1 %
28	Lead sulfochromate yellow (C.I. Pigment Yellow 34)	1344-37-2	215-693-7	< 0.1 %
29	Ammonium dichromate	7789-09-5	232-143-1	< 0.1 %
30	Boric acid	10043-35-3 11113-50-1	233-139-2 234-343-4	< 0.1 %
31	Disodium tetraborate, anhydrous	1303-96-4 1330-43-4 12179-04-3	215-540-4	< 0.1 %
32	Potassium chromate	7789-00-6	232-140-5	< 0.1 %
33	Potassium dichromate	7778-50-9	231-906-6	< 0.1 %
34	Sodium chromate	7775-11-3	231-889-5	< 0.1 %
35	Tetraboron disodium heptaoxide, hydrate	12267-73-1	235-541-3	< 0.1 %
36	Trichloroethylene	79-01-6	201-167-4	< 0.1 %
37	Cobalt(II) sulphate	10124-43-3	233-334-2	< 0.1 %
38	Cobalt(II) dinitrate	10141-05-6	233-402-1	< 0.1 %
39	Cobalt(II) carbonate	513-79-1	208-169-4	< 0.1 %
40	Cobalt(II) diacetate	71-48-7	200-755-8	< 0.1 %
41	2-Methoxyethanol	109-86-4	203-713-7	< 0.1 %
42	2-Ethoxyethanol	110-80-5	203-804-1	< 0.1 %
43	Chromium trioxide	1333-82-0	215-607-8	< 0.1 %
44	Acids generated from chromium trioxide and their oligomers: Chromic acid Dichromic acid Oligomers of chromic acid and dichromic acid	7738-94-5 13530-68-2 -	231-801-5 236-881-5 -	< 0.1 %
45	2-ethoxyethyl acetate	111-15-9	203-839-2	< 0.1 %
46	Strontium chromate	7789-06-2	232-142-6	< 0.1 %
47	1,2-Benzenedicarboxylic acid, di-C7-11-branched and li near alkyl esters (DNHUP)	68515-42-4	271-084-6	< 0.1 %
48	Hydrazine	7803-57-8 302-01-2	206-114-9	< 0.1 %
49	1-methyl-2-pyrrolidone	872-50-4	212-828-1	< 0.1 %
50	1,2,3-trichloropropane	96-18-4	202-486-1	< 0.1 %
51	1,2-Benzenedicarboxylic acid, di-C6-8-branched alkyl esters, C7-rich	71888-89-6	276-158-1	< 0.1 %
52	Lead dipicrate	6477-64-1	229-335-2	< 0.1 %
53	Lead styphnate	15245-44-0	239-290-0	< 0.1 %
54	Lead diazide, Lead azide	13424-46-9	236-542-1	< 0.1 %
55	Phenolphthalein	77-09-8	201-004-7	< 0.1 %
56	2,2'-dichloro-4,4'-methylenedianiline (MOCA)	101-14-4	202-918-9	< 0.1 %
57	N,N-dimethylacetamide (DMAC)	127-19-5	204-826-4	< 0.1 %

58	Trilead diarsenate	3687-31-8	222-979-5	< 0.1 %
59	Calcium arsenate	7778-44-1	231-904-5	< 0.1 %
60	Arsenic acid	7778-39-4	231-901-9	< 0.1 %
61	Bis(2-methoxyethyl) ether	111-96-6	203-924-4	< 0.1 %
62	1,2-Dichloroethane	107-06-2	203-458-1	< 0.1 %
63	4-(1,1,3,3-tetramethylbutyl)phenol	140-66-9	205-426-2	< 0.1 %
64	2-Methoxyaniline; o-Anisidine	90-04-0	201-963-1	< 0.1 %
65	Bis(2-methoxyethyl) phthalate	117-82-8	204-212-6	< 0.1 %
66	Formaldehyde, oligomeric reaction products with aniline (technical MDA)	25214-70-4	500-036-1	< 0.1 %
67	Aluminosilicate Refractory Ceramic Fibres	-	-	< 0.1 %
68	Zirconia Aluminosilicate Refractory Ceramic Fibres	-	-	< 0.1 %
69	Pentazinc chromate octahydroxide	49663-84-5	256-418-0	< 0.1 %
70	Potassium hydroxyoctaoxodizincatedichromate	11103-86-9	234-329-8	< 0.1 %
71	Dichromium tris(chromate)	24613-89-6	246-356-2	< 0.1 %
72	1,2-bis(2-methoxyethoxy)ethane (TEGDME; triglyme)	112-49-2	203-977-3	< 0.1 %
73	1,2-dimethoxyethane; ethylene glycol dimethyl ether (EGDME)	110-71-4	203-794-9	< 0.1 %
74	Diboron trioxide	1303-86-2	215-125-8	< 0.1 %
75	Formamide	75-12-7	200-842-0	< 0.1 %
76	Lead(II) bis(methanesulfonate)	17570-76-2	401-750-5	< 0.1 %
77	TGIC (1,3,5-tris(oxiranylmethyl)-1,3,5-triazine-2,4,6(1H,3H,5H)-trione)	2451-62-9	219-514-3	< 0.1 %
78	β-TGIC (1,3,5-tris[(2S and 2R)-2,3-epoxypropyl]-1,3,5-triazine-2,4,6-(1H,3H,5H)-trione)	59653-74-6	423-400-0	< 0.1 %
79	4,4'-bis(dimethylamino)benzophenone (Michler's ketone)	90-94-8	202-027-5	< 0.1 %
80	N,N,N',N'-tetramethyl-4,4'-methylenedianiline (Michler's base)	101-61-1	202-959-2	< 0.1 %
81	[4-[4,4'-bis(dimethylamino) benzhydrylidene]cyclohexa-2,5-dien-1-ylidene]dimethylammonium chloride (C.I. Basic Violet 3)	548-62-9	208-953-6	< 0.1 %
82	[4-[[4-anilino-1-naphthyl][4-(dimethylamino)phenyl]methylene]cyclohexa-2,5-dien-1-ylidene] dimethylammonium chloride (C.I. Basic Blue 26)	2580-56-5	219-943-6	< 0.1 %
83	α,α-Bis[4-(dimethylamino)phenyl]-4 (phenylamino)naphthalene-1-methanol (C.I. Solvent Blue 4)	6786-83-0	229-851-8	< 0.1 %
84	4,4'-bis(dimethylamino)-4''-(methylamino)trityl alcohol	561-41-1	209-218-2	< 0.1 %
85	Pyrochlore, antimony lead yellow	8012-00-8	232-382-1	< 0.1 %
86	6-methoxy-m-toluidine (p-cresidine)	120-71-8	204-419-1	< 0.1 %
87	Henicosafluoroundecanoic acid	2058-94-8	218-165-4	< 0.1 %
88	Hexahydromethylphthalic anhydride [1], Hexahydro-4-methylphthalic anhydride [2], Hexahydro-1-methylphthalic anhydride [3],	25550-51-0 19438-60-9 48122-14-1	247-094-1 243-072-0 256-356-4	< 0.1 %

	Hexahydro-3-methylphthalic anhydride [4] [The individual isomers [2], [3] and [4] (including their cis- and trans- stereo isomeric forms) and all possible combinations of the isomers [1] are covered by this entry]	57110-29-9	260-566-1	
89	Cyclohexane-1,2-dicarboxylic anhydride [1], cis-cyclohexane-1,2-dicarboxylic anhydride [2], trans-cyclohexane-1,2-dicarboxylic anhydride [3] [The individual cis- [2] and trans- [3] isomer substances and all possible combinations of the cis- and trans-isomers [1] are covered by this entry]	85-42-7 13149-00-3 14166-21-3	201-604-9 236-086-3 238-009-9	< 0.1 %
90	Dibutyltin dichloride (DBTC)	683-18-1	211-670-0	< 0.1 %
91	Lead bis(tetrafluoroborate)	13814-96-5	237-486-0	< 0.1 %
92	Lead dinitrate	10099-74-8	233-245-9	< 0.1 %
93	Silicic acid, lead salt	11120-22-2	234-363-3	< 0.1 %
94	4-Aminoazobenzene	60-09-3	200-453-6	< 0.1 %
95	Lead titanium zirconium oxide	12626-81-2	235-727-4	< 0.1 %
96	Lead monoxide (lead oxide)	1317-36-8	215-267-0	< 0.1 %
97	o-Toluidine	95-53-4	202-429-0	< 0.1 %
98	3-ethyl-2-methyl-2-(3-methylbutyl)-1,3-oxazolidine	143860-04-2	421-150-7	< 0.1 %
99	Silicic acid (H2Si2O5), barium salt (1:1), lead-doped [with lead (Pb) content above the applicable generic concentration limit for 'toxicity for reproduction' Repr. 1A (CLP) or category 1 (DSD); the substance is a member of the group entry of lead compounds, with index number 082-001-00-6 in Regulation (EC) No 1272/2008]	68784-75-8	272-271-5	< 0.1 %
100	Trilead bis(carbonate)dihydroxide	1319-46-6	215-290-6	< 0.1 %
101	Furan	110-00-9	203-727-3	< 0.1 %
102	N,N-dimethylformamide	68-12-2	200-679-5	< 0.1 %
103	4-(1,1,3,3-tetramethylbutyl)phenol, ethoxylated [covering well-defined substances and UVCB substances, polymers and homologues]	-	-	< 0.1 %
104	4-Nonylphenol, branched and linear [substances with a linear and/or branched alkyl chain with a carbon number of 9 covalently bound in position 4 to phenol, covering also UVCB- and well-defined substances which include any of the individual isomers or a combination thereof]	-	-	< 0.1 %
105	4,4'-methylenedi-o-toluidine	838-88-0	212-658-8	< 0.1 %
106	Diethyl sulphate	64-67-5	200-589-6	< 0.1 %
107	Dimethyl sulphate	77-78-1	201-058-1	< 0.1 %
108	Lead oxide sulfate	12036-76-9	234-853-7	< 0.1 %
109	Lead titanium trioxide	12060-00-3	235-038-9	< 0.1 %
110	Acetic acid, lead salt, basic	51404-69-4	257-175-3	< 0.1 %
111	[Phthalato(2-)]dioxotrilead	69011-06-9	273-688-5	< 0.1 %
112	Bis(pentabromophenyl) ether (decabromodiphenyl ether; DecaBDE)	1163-19-5	214-604-9	< 0.1 %

113	N-methylacetamide	79-16-3	201-182-6	< 0.1 %
114	Dinoseb (6-sec-butyl-2,4-dinitrophenol)	88-85-7	201-861-7	< 0.1 %
115	1,2-Diethoxyethane	629-14-1	211-076-1	< 0.1 %
116	Tetralead trioxide sulphate	12202-17-4	235-380-9	< 0.1 %
117	N-pentyl-isopentylphthalate	776297-69-9	-	< 0.1 %
118	Dioxobis(stearato)trilead	12578-12-0	235-702-8	< 0.1 %
119	Tetraethyllead	78-00-2	201-075-4	< 0.1 %
120	Pentalead tetraoxide sulphate	12065-90-6	235-067-7	< 0.1 %
121	Pentacosafluorotridecanoic acid	72629-94-8	276-745-2	< 0.1 %
122	Tricosafluorododecanoic acid	307-55-1	206-203-2	< 0.1 %
123	Heptacosafluorotetradecanoic acid	376-06-7	206-803-4	< 0.1 %
124	1-bromopropane (n-propyl bromide)	106-94-5	203-445-0	< 0.1 %
125	Methoxyacetic acid	625-45-6	210-894-6	< 0.1 %
126	4-methyl-m-phenylenediamine (toluene-2,4-diamine)	95-80-7	202-453-1	< 0.1 %
127	Methyloxirane (Propylene oxide)	75-56-9	200-879-2	< 0.1 %
128	Trilead dioxide phosphonate	12141-20-7	235-252-2	< 0.1 %
129	o-aminoazotoluene	97-56-3	202-591-2	< 0.1 %
130	1,2-Benzenedicarboxylic acid, dipentylester, branched and linear	84777-06-0	284-032-2	< 0.1 %
131	4,4'-oxydianiline and its salts	101-80-4	202-977-0	< 0.1 %
132	Orange lead (lead tetroxide)	1314-41-6	215-235-6	< 0.1 %
133	Biphenyl-4-ylamine	92-67-1	202-177-1	< 0.1 %
134	Diisopentylphthalate	605-50-5	210-088-4	< 0.1 %
135	Fatty acids, C16-18, lead salts	91031-62-8	292-966-7	< 0.1 %
136	Diazene-1,2-dicarboxamide (C,C'-azodi(formamide))	123-77-3	204-650-8	< 0.1 %
137	Sulfurous acid, lead salt, dibasic	62229-08-7	263-467-1	< 0.1 %
138	Lead cyanamidate	20837-86-9	244-073-9	< 0.1 %
139	Cadmium	7440-43-9	231-152-8	< 0.1 %
140	Ammonium pentadecafluorooctanoate (APFO)	3825-26-1	223-320-4	< 0.1 %
141	Pentadecafluorooctanoic acid (PFOA)	335-67-1	206-397-9	< 0.1 %
142	Dipentyl phthalate (DPP)	131-18-0	205-017-9	< 0.1 %
143	4-Nonylphenol, branched and linear, ethoxylated [substances with a linear and/or branched alkyl chain with a carbon number of 9 covalently bound in position 4 to phenol, ethoxylated covering UVCB- and well-defined substances, polymers and homologues, which include any of the individual isomers and/or combinations thereof]	-	-	< 0.1 %
144	Cadmium oxide	1306-19-0	215-146-2	< 0.1 %
145	Cadmium sulphide	1306-23-6	215-147-8	< 0.1 %
146	Disodium 4-amino-3-[[4'-[(2,4-diaminophenyl)azo][1,1'-biphenyl]-4-yl]azo] -5-hydroxy-6-(phenylazo)naphthalene-2,7-disulphonate	1937-37-7	217-710-3	< 0.1 %

	(C.I. Direct Black 38)			
147	Dihexyl phthalate	84-75-3	201-559-5	< 0.1 %
148	Imidazolidine-2-thione; (2-imidazoline-2-thiol)	96-45-7	202-506-9	< 0.1 %
149	Trixylyl phosphate	25155-23-1	246-677-8	< 0.1 %
150	Disodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis(azo)]bis(4-aminonaphthalene-1-sulphonate) (C.I. Direct Red 28)	573-58-0	209-358-4	< 0.1 %
151	Lead di(acetate)	301-04-2	206-104-4	< 0.1 %
152	1,2-Benzenedicarboxylic acid, dihexyl ester, branched and linear	68515-50-4	271-093-5	< 0.1 %
153	Sodium perborate; perboric acid, sodium salt	-	239-172-9 234-390-0	< 0.1 %
154	Sodium peroxometaborate	7632-04-4	231-556-4	< 0.1 %
155	Cadmium chloride	10108-64-2	233-296-7	< 0.1 %
156	2-benzotriazol-2-yl-4,6-di-tert-butylphenol (UV-320)	3846-71-7	223-346-6	< 0.1 %
157	2-ethylhexyl 10-ethyl-4, 4-dioctyl-7-oxo-8-oxa-3, 5-dithia-4-stannatetradecanoate (DOTE)	15571-58-1	239-622-4	< 0.1 %
158	Reaction mass of 2-ethylhexyl 10-ethyl-4,4-dioctyl-7-oxo-8-oxa-3,5-dithia-4-stannatetradecanoate and 2-ethylhexyl 10-ethyl-4-[[2-[(2-ethylhexyl)oxy]-2-oxoethyl]thio]-4-octyl-7-oxo-8-oxa-3,5-dithia-4-stannatetradecanoate (reaction mass of DOTE and MOTE)	-	-	< 0.1 %
159	2-(2H-benzotriazol-2-yl)-4,6-ditertpentylphenol (UV-328)	25973-55-1	247-384-8	< 0.1 %
160	Cadmium fluoride	7790-79-6	232-222-0	< 0.1 %
161	Cadmium sulphate	10124-36-4 31119-53-6	233-331-6	< 0.1 %
162	Reaction mass of 2-ethylhexyl 10-ethyl-4,4-dioctyl-7-oxo-8-oxa-3,5-dithia-4-stannatetradecanoate and 2-ethylhexyl 10-ethyl-4-[[2-[(2-ethylhexyl)oxy]-2-oxoethyl]thio]-4-octyl-7-oxo-8-oxa-3,5-dithia-4-stannatetradecanoate (reaction mass of DOTE and MOTE)	68515-51-5 68648-93-1	271-094-0 272-013-1	< 0.1 %
163	5-sec-butyl-2-(2,4-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane [1], 5-sec-butyl-2-(4,6-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane [2] [covering any of the individual stereoisomers of [1] and [2] or any combination thereof]	_	_	< 0.1 %
164	Nitrobenzene	98-95-3	202-716-0	< 0.1 %
165	2,4-di-tert-butyl-6-(5-chlorobenzotriazol-2-yl)phenol (UV-327)	3864-99-1	223-383-8	< 0.1 %
166	2-(2H-benzotriazol-2-yl)-4-(tert-butyl)-6-(sec-butyl)phenol (UV-350)	36437-37-3	253-037-1	< 0.1 %
167	1,3-propanesultone	1120-71-4	214-317-9	< 0.1 %
168	Perfluorononan-1-oic-acid and its sodium and ammonium salts	375-95-1 21049-39-8 4149-60-4	206-801-3	< 0.1 %

169	Benzo [def] chrysene	50-32-8	200-028-5	< 0.1 %
170	4,4'-isopropylidenediphenol Bisphenol A; BPA	80-05-7	201-245-8	< 0.1 %
171	4-heptylphenol, branched and linear substances with a linear and/or branched alkyl chain with a carbon number of 7 covalently bound predominantly in position 4 to phenol, covering also UVCB- and well-defined substances which include any of the individual isomers or a combination thereof	-	-	< 0.1 %
172	Nonadecafluorodecanoic acid (PFDA) and its sodium and ammonium salts	3108-42-7 335-76-2 3830-45-3	206-400-3 221-470-5	< 0.1 %
173	p-(1,1-dimethylpropyl)phenol	80-46-6	201-280-9	< 0.1 %
174	Perfluorohexane-1-sulphonic acid and its salts PFHxS	-	-	< 0.1 %
175	1,6,7,8,9,14,15,16,17,17,18,18-Dodecachloropentacyclo[12.2.1.16,9.02,13.05,10]octadeca-7,15-diene (“Dechlorane Plus”™) covering any of its individual anti- and syn-isomers or any combination thereof	-	-	< 0.1 %
176	Benz[a]anthracene	56-55-3 1718-53-2	200-280-6	< 0.1 %
177	Cadmium carbonate	56-55-3 1718-53-2	200-280-6	< 0.1 %
178	Cadmium hydroxide	21041-95-2	244-168-5	< 0.1 %
179	Cadmium nitrate	10022-68-1 10325-94-7	233-710-6	< 0.1 %
180	Chrysene	218-01-9 1719-03-5	205-923-4	< 0.1 %
181	Reaction products of 1,3,4-thiadiazolidine-2,5-dithione, formaldehyde and 4-heptylphenol, branched and linear (RP-HP) with ≥0.1% w/w 4-heptylphenol, branched and linear (4-HPbl)	-	-	< 0.1 %
182	Benzene-1,2,4-tricarboxylic acid 1,2 anhydride	552-30-7	209-008-0	< 0.1 %
183	Benzo[ghi]perylene	191-24-2	205-883-8	< 0.1 %
184	Decamethylcyclopentasiloxane	541-02-6	208-764-9	< 0.1 %
185	Dicyclohexyl phthalate	84-61-7	201-545-9	< 0.1 %
186	Disodium octaborate	12008-41-2	234-541-0	< 0.1 %
187	Dodecamethylcyclohexasiloxane	540-97-6	208-762-8	< 0.1 %
188	Ethylenediamine	107-15-3	203-468-6	< 0.1 %
189	Lead	7439-92-1	231-100-4	< 0.1 %
190	Octamethylcyclotetrasiloxane	556-67-2	209-136-7	< 0.1 %
191	Terphenyl, hydrogenated	61788-32-7	262-967-7	< 0.1 %
192	Pyrene	129-00-0 1718-52-1	204-927-3	< 0.1 %
193	Phenanthrene	85-01-8	201-581-5	< 0.1 %
194	Fluoranthene	206-44-0 93951-69-0	205-912-4	< 0.1 %
195	Benzo [k] fluoranthene	207-08-9	205-916-6	< 0.1 %

196	2,2-bis(4'-hydroxyphenyl)-4-methylpentane	6807-17-6	401-720-1	< 0.1 %
197	1,7,7-trimethyl-3-(phenylmethylene) bicycle [2.2.1] heptan-2-one	15087-24-8	239-139-9	< 0.1 %
198	2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)propionic acid, its salts and its acyl halides	-	-	< 0.1 %
199	2-methoxyethyl acetate	110-49-6	203-772-9	< 0.1 %
200	4-tert-butylphenol	98-54-4	202-679-0	< 0.1 %
201	Tris(4-nonylphenyl, branched and linear) phosphite (TNPP) with ≥ 0.1% w/w of 4-nonylphenol, branched and linear (4-NP)	-	-	< 0.1 %
202	2-benzyl-2-dimethylamino-4'-morpholinobutyrophenone	119313-12-1	404-360-3	< 0.1 %
203	2-methyl-1-(4-methylthiophenyl)-2-morpholinopropan-1-one	71868-10-5	400-600-6	< 0.1 %
204	Diisohexyl phthalate	71850-09-4	276-090-2	< 0.1 %
205	Perfluorobutane sulfonic acid (PFBS) and its salts	-	-	< 0.1 %
206	1-vinylimidazole	1072-63-5	214-012-0	< 0.1 %
207	2-methylimidazole	693-98-1	211-765-7	< 0.1 %
208	Butyl 4-hydroxybenzoate	94-26-8	202-318-7	< 0.1 %
209	Dibutylbis(pentane-2,4-dionato-O,O')tin	22673-19-4	245-152-0	< 0.1 %
210	Bis(2-(2-methoxyethoxy)ethyl) ether	143-24-8	205-594-7	< 0.1 %
211	Dioctyltin dilaurate, stannane, dioctyl-, bis(coco acyloxy) derivs., and any other stannane, dioctyl-, bis(fatty acyloxy) derivs. wherein C12 is the predominant carbon number of the fatty acyloxy moiety ・dioctyltin dilaurate; stannane, dioctyl-, bis(coco acyloxy) derivs. ・Stannane, dioctyl-, bis(coco acyloxy) derivs. ・Dioctyltin dilaurate	- 91648-39-4 3648-18-8	-	< 0.1 %
212	1,4-dioxane	123-91-1	204-661-8	< 0.1 %
213	2,2-bis(bromomethyl)propane-1,3-diol (BMP); 2,2-dimethylpropan-1-ol, tribromo derivative/3-bromo-2,2-bis(bromomethyl)-1-propanol (TBNPA); 2,3-dibromo-1-propanol (2,3-DBPA)	3296-90-0 36483-57-5/1522-92-5 96-13-9	221-967-7 253-057-0/- 202-480-9	< 0.1 %
214	2-(4-tertbutylbenzyl)propionaldehyde and its individual stereoisomers	75166-31-3 80-54-6 75166-30-2	- 201-289-8 -	< 0.1 %
215	4,4'-(1-methylpropylidene)bisphenol	77-40-7	201-025-1	< 0.1 %
216	glutaral	111-30-8	203-856-5	< 0.1 %
217	Medium-chain chlorinated paraffins(MCCP) [UVCB substances consisting of more than or equal to 80% linear chloroalkanes with carbon chain lengths within the range from C14 to C17]	1372804-76-6 85535-85-9 - 198840-65-2	- 287-477-0 950-299-5 -	< 0.1 %
218	orthoboric acid, sodium salt	25747-83-5 22454-04-2 14312-40-4 1333-73-9 13840-56-7 14890-53-0	- - 238-253-6 215-604-1 237-560-2 -	< 0.1 %
219	Phenol, alkylation products (mainly in para position) with	210555-94-5	-	< 0.1 %

	C12-rich branched alkyl chains from oligomerisation, covering any individual isomers and/ or combinations thereof (PDDP)	27459-10-5 27147-75-7 121158-58-5 74499-35-7 57427-55-	- - 310-154-3 - -	
220	(±)-1,7,7-trimethyl-3-[(4-methylphenyl)methylene]bicyclo[2.2.1] heptan-2-one covering any of the individual isomers and/or combinations thereof (4-MBC)	1782069-81-1 95342-41-9 852541-25-4 36861-47-9 741687-98-9 852541-30-1 852541-21-0	- - - 253-242-6 - - -	< 0.1 %
221	6,6'-di-tert-butyl-2,2'-methylenedi-p-cresol	119-47-1	204-327-1	< 0.1 %
222	S-(tricyclo(5.2.1.0'2,6)deca-3-en-8(or 9)-yl O-(isopropyl or isobutyl or 2-ethylhexyl) O-(isopropyl or isobutyl or 2-ethylhexyl) phosphorodithioate	255881-94-8	401-850-9	< 0.1 %
223	tris(2-methoxyethoxy)vinylsilane	1067-53-4	213-934-0	< 0.1 %
224	N-(hydroxymethyl)acrylamide	924-42-5	213-103-2	< 0.1 %

Signature: Alan Shen

Alan Shen
Director of QA Division

Date: July 28, 2022

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