Règlement (UE) no 1907/2006 - REACH

REACH est un réglement de l'Union Européenne sur les substances chimiques, leur maintient sous contrôle et leur sécurité d'utilisation. Il concerne l'enregistrement, l'évaluation, l'autorisation et la restriction des substances chimiques (REACH est l'acronime de **R**egistration, **E**valuation, **A**uthorisation and Restriction of **Ch**emicals).

REACH remplace de nombreux textes législatifs de l'Union relatifs aux substances chimiques (ainsi que les composants contenant ces substances) et complète d'autres textes concernant l'environnement et la sécurité.

L'un des objectifs les plus importants est sûrement l'amélioration de la gestion de la santé humaine et de l'environnement face aux risques potentiels liés à l'utilisation des substances chimiques elles-mêmes, ou des composants d'un mélange ou d'un article.

Ces restrictions se situent dans un contexte intégré par Finder depuis longtemps déjà. Comme **utilisateurs finaux (en aval)** de substances chimiques, nous communiquons activement avec nos fournisseurs en analysant et en tenant sous étroit contrôle les fiches de sécurité prévues pour chaque type desubstance ou de préparation. À noter que Finder emploie des substances et/ou des préparations dans ses processus de fabrication sans en modifier les caractéristiques chimiques.

Dans les délais d'adaptation aux obligations prévues par la Directive, Finder s'engage à maintenir une grande attention sur le choix de ses fournisseurs et garantit une communication optimale avec ses propres Clients en se référant scrupuleusement aux mises à jour des tableaux publiés par les organismes officiels (ECHA) – En pièce jointe la liste des substances hautement dangereuses "Candidate list of substances of very high concern for authorization (SVHC)" et l'inventaire des classifications des substances du document joint XVII du Règlement Reach mis à jour.

En outre, cette politique est mise en œuvre par un Système ISO 14001 (Système de Gestion Environnementale), qui utilise des questionnaires destinés à évaluer préventivement les personnes qui collaborent avec Finder. En tant qu'**utilisateurs aval** de substances chimiques, elles doivent prendre des mesures préventives contre les risques identifiés. Pour ce motif, en plus de ce qui est spécifié plus haut, l'entreprise a mis en place un Système de Gestion Santé et Sécurité qui impose une prise en compte particulièrement sensible des questions et initiatives dans ce domaine afin de garantir la protection des personnes impliquées dans les divers processus de fabrication.

REACH – Substances extrêmement dangereuses
Candidate List of Substances of Very High Concern – SVHC

https://echa.europa.eu/candidate-list-table

FINDER précise que ses produits ne contiennent aucune des substances indiquées dans le tableau reproduit ci-dessous SVHC en concentration supérieure à 0.1% en poids.

Noms des Substances	Numéro EC	Numéro CAS	Numéro EC	Numéro CAS
Diphenyl(2,4,6-trimethylbenzoyl)phosphine oxide			278-355-8	75980-60-8
Bis(4-chlorophenyl) sulphone			201-247-9	80-07-9
reaction mass of 2,2,3,3,5,5,6,6-octafluoro-4-(1,1,1,2,3,3,3-heptafluoropropan-2-yl)morpholine and 2,2,3,3,5,5,6,6-octafluoro-4-(heptafluoropropyl)morpholine			473-390-7	-
Perfluoroheptanoic acid and its salts			-	-
Ammonium perfluoroheptanoate	228-098-2	6130-43-4		
potassium perfluoroheptanoate	-	21049-36-5		
Perfluoroheptanoic acid	206-798-9	375-85-9		
Sodium perfluoroheptanoate	243-518-4	20109-59-5		
Melamine			203-615-4	108-78-1
Isobutyl 4-hydroxybenzoate			224-208-8	4247-02-3
bis(2-ethylhexyl) tetrabromophthalate covering any of the individual isomers and/or combinations thereof			-	-
Bis(2-ethylhexyl) tetrabromophthalate	247-426-5	26040-51-7		
Barium diboron tetraoxide			237-222-4	13701-59-2
4,4'-sulphonyldiphenol			201-250-5	80-09-1
2,2',6,6'-tetrabromo-4,4'-isopropylidenediphenol			201-236-9	79-94-7
1,1'-[ethane-1,2-diylbisoxy]bis[2,4,6-tribromobenzene]			253-692-3	37853-59-1
N-(hydroxymethyl)acrylamide			213-103-2	924-42-5
tris(2-methoxyethoxy)vinylsilane				
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 X4261			924-42-5	255881-94-8
6,6'-di-tert-butyl-2,2'-methylenedi-p-cresol			204-327-1	119-47-1
(±)-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)			-	-
(3E)-1,7,7-trimethyl-3-(4-methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	1782069-81-1		
(1R,3E,4S)-1,7,7-trimethyl-3-(4- methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	95342-41-9		
(1S,3Z,4R)-1,7,7-trimethyl-3-(4-methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	852541-25-4		
(±)-1,7,7-trimethyl-3-[(4-methylphenyl)methylene] bicyclo[2.2.1]heptan-2-one	253-242-6	36861-47-9		
(1R,4S)-1,7,7-trimethyl-3-(4-methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	741687-98-9		
(1S,3E,4R)-1,7,7-trimethyl-3-(4- methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	852541-30-1		
(1R,3Z,4S)-1,7,7-trimethyl-3-(4- methylbenzylidene) bicyclo[2.2.1]heptan-2-one	-	852541-21-0		
Phenol, alkylation products (mainly in para position) with C12-rich branched alkyl chains from oligomerisation, covering any individual isomers and/ or combinations thereof (PDDP)			-	-
Phenol, 4-dodecyl, branched	-	210555-94-5		
4-isododecylphenol	-	27459-10-5		
Phenol, 4-isododecyl-	-	27147-75-7		
Phenol, dodecyl-, branched	310-154-3	121158-58-5		
Phenol, (tetrapropenyl) derivatives	-	74499-35-7		
Phenol, tetrapropylene-	-	57427-55-1		
orthoboric acid, sodium salt			-	-
boric acid (H3BO3), sodium salt, hydrate	-	25747-83-5		
Boric acid (H3BO3), disodium salt	-	22454-04-2		

Trisodium orthoborate	238-253-6	14312-40-4		
Boric acid, sodium salt	215-604-1	1333-73-9		
Orthoboric acid, sodium salt	237-560-2	13840-56-7		
Boric acid (H3BO3), sodium salt (1:1)	-	14890-53-0		
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			-	-
Alkanes, C14-16, chloro	-	1372804-76-6		
Alkanes, C14-17, chloro	287-477-0	85535-85-9		
di-, tri- and tetrachlorotetradecane	950-299-5	-		
Tetradecane, chloro derivs.	-	198840-65-2		
glutaral			203-856-5	111-30-8
4,4'-(1-methylpropylidene)bisphenol			201-025-1	77-40-7
2-(4-tert-butylbenzyl)propionaldehyde and its individual stereoisomers			-	-
(2R)-3-(4-tert-butylphenyl)-2-methylpropanal	-	75166-31-3		
2-(4-tert-butylbenzyl)propionaldehyde	201-289-8	80-54-6		
(2S)-3-(4-tert-butylphenyl)-2-methylpropanal	-	75166-30-2		
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)			-	-
2,2-dimethylpropan-1-ol, tribromo derivative (TBNPA)	253-057-0	36483-57-5		
3-bromo-2,2-bis(bromomethyl)-1-propanol (TBNPA)	-	1522-92-5		
2,2-bis(bromomethyl)propane-1,3-diol (BMP)	221-967-7	3296-90-0		
2,3-dibromo-1-propanol (2,3-DBPA)	202-480-9	96-13-9		
1,4-dioxane			204-661-8	123-91-1
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			-	-
Stannane, dioctyl-, bis(coco acyloxy) derivs.	293-901-5	91648-39-4		
dioctyltin dilaurate; stannane, dioctyl-, bis(coco acyloxy) derivs.	-	-		
Dioctyltin dilaurate	222-883-3	3648-18-8		
Bis(2-(2-methoxyethoxy)ethyl)ether			205-594-7	143-24-8
Dibutylbis(pentane-2,4-dionato-O,O')tin			245-152-0	22673-19-4
Butyl 4-hydroxybenzoate			202-318-7	94-26-8
2-methylimidazole			211-765-7	693-98-1
1-vinylimidazole			214-012-0	1072-63-5
Perfluorobutane sulfonic acid (PFBS) and its salts			-	-
bis(4-t-butylphenyl)iodonium perfluorobutanesulfonate	432-660-4	-		
tetrabutyl-phosphonium nonafluoro-butane-1-sulfonate	444-440-5	220689-12-3		
dimethyl(phenyl)sulfanium perfluorobutanesulfonate	452-310-4	220133-51-7		
1-(4-butoxy-1-naphthalenyl)tetrahydrothiophenium 1,1,2,2,3,3,4,4,4-nonafluoro-1-butanesulfonate	468-770-4	-		
Triphenylsulfanium perfluorobutane sulfonate	478-340-8	144317-44-2		
1,1,2,2,3,3,4,4,4-nonafluorobutane-1-sulphonic acid	206-793-1	375-73-5		
lithium perfluorobutanesulfonate	-	131651-65-5		
morpholinium perfluorobutanesulfonate	-	503155-89-3		
N,N,N-triethylethanaminium 1,1,2,2,3,3,4,4,4-nonafluorobutane-1-sulfonate	-	25628-08-4		
Potassium 1,1,2,2,3,3,4,4,4-nonafluorobutane-1-sulphonate	249-616-3	29420-49-3		
Ammonium 1,1,2,2,3,3,4,4,4-nonafluorobutane-1-sulphonate	269-513-7	68259-10-9		
magnesium perfluorobutanesulfonate	-	507453-86-3		
Diisohexyl phthalate			276-090-2	71850-09-4
2-methyl-1-(4-methylthiophenyl)-2-morpholinopropan-1-one			400-600-6	71868-10-5
2-benzyl-2-dimethylamino-4'-morpholinobutyrophenone			404-360-3	119313-12-1
Tris(4-nonylphenyl, branched and linear) phosphite (TNPP) with ≥ 0.1% w/w of 4-nonylphenol, branched and linear (4-NP)			-	-

Phenol, 4-nonyl-, phosphite (3:1)	-	3050-88-2		
tris(4-nonylphenyl, branched) phosphite	-	-		
tris(nonylphenyl) phosphite	247-759-6	26523-78-4		
Phenol, p-isononyl-, phosphite (3:1)	-	31631-13-7		
Phenol, p-sec-nonyl-, phosphite	-	106599-06-8		
4-tert-butylphenol			202-679-0	98-54-4
2-methoxyethyl acetate			203-772-9	110-49-6
2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)propionic acid, its salts and its acyl halides covering any of their individual isomers and combinations thereof			-	-
potassium 2,3,3,3-tetrafluoro-2-(heptafluoropropoxy) propionate	266-578-3	67118-55-2		
2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)propionyl fluoride	218-173-8	2062-98-8		
ammonium 2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)propanoate	-	62037-80-3		
2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)propionic acid	236-236-8	13252-13-6		
Propanoic acid, 2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)-, (-)-	-	75579-40-7		
Propanoic acid, 2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)-, (+)-	-	75579-39-4		
Pyrene			204-927-3	129-00-0
Phenanthrene			201-581-5	85-01-8
Fluoranthene			205-912-4	206-44-0
Benzo[k]fluoranthene			205-916-6	207-08-9
2,2-bis(4'-hydroxyphenyl)-4-methylpentane			401-720-1	6807-17-6
1,7,7-trimethyl-3-(phenylmethylene)bicyclo[2.2.1]heptan-2-one 3-benzylidene camphor; 3-BC			239-139-9	15087-24-8
Terphenyl, hydrogenated			262-967-7	61788-32-7
Octamethylcyclotetrasiloxane D4			209-136-7	556-67-2
Lead			231-100-4	7439-92-1
Ethylenediamine EDA			203-468-6	107-15-3
Dodecamethylcyclohexasiloxane D6			208-762-8	540-97-6
Disodium octaborate			234-541-0	12008-41-2
Dicyclohexyl phthalate DCHP			201-545-9	84-61-7
Decamethylcyclopentasiloxane D5			208-764-9	541-02-6
Benzo[ghi]perylene			205-883-8	191-24-2
Benzene-1,2,4-tricarboxylic acid 1,2 anhydride trimellitic anhydride; TMA			209-008-0	552-30-7
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)			-	-
Formaldehyde, reaction products with branched and linear heptylphenol, carbon disulfide and hydrazine	300-298-5	93925-00-9		
Reaction product of 1,3,4-thiadiazolidine-2,5-dithione, formaldehyde and phenol, heptyl derivs.	-	1471311-26-8		
Chrysene			205-923-4	218-01-9
Cadmium nitrate			233-710-6	10325-94-7
Cadmium hydroxide			244-168-5	21041-95-2
Cadmium carbonate			208-168-9	513-78-0
Benz[a]anthracene			200-280-6	56-55-3
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			-	-
1,6,7,8,9,14,15,16,17,17,18,18-dodecachloropen-tacyclo[12.2.1.16,9.02,13.05,10]octadeca-7,15-diene	236-948-9	13560-89-9		

(1S,2S,5S,6S,9R,10R,13R,14R)-1,6,7,8,9,14,15,16,17,17,18,18-Dodecachloropentacyclo[12.2.1.1^{6,9}.0^{2,13}.0^{5,10}]octadeca-7,15-diene	-	135821-74-8		
(1S,2S,5R,6R,9S,10S,13R,14R)-1,6,7,8,9,14,15,16,17,17,18,18-Dodecachloropentacyclo[12.2.1.1^{6,9}.0^{2,13}.0^{5,10}]octadeca-7,15-diene	-	135821-03-3		
rel-(1R,4S,4aS,6aS,7S,10R,10aR,12aR)-1,2,3,4,7,8,9,10,-13,13,14,14-dodecachloro-1,4,4a,5,6,6a,7,10,10a,11,12,12a-dodecahydro-1,4:7,10-dimethanodibenzo[a,e]cyclooctene	-	-		
rel-(1R,4S,4aS,6aR,7R,10S,10aS,12aR)-1,2,3,4,7,8,9,10,-13,13,14,14-dodecachloro-1,4,4a,5,6,6a,7,10,10a,11,12,12a-dodecahydro-1,4:7,10-dimethanodibenzo[a,e]cyclooctene	-	-		
Perfluorohexane-1-sulphonic acid and its salts PFHxS			-	-
tridecafluorohexanesulphonic acid, compound with 2,2'-iminodiethanol (1:1)	274-462-9	70225-16-0		
ammonium perfluorohexane-1-sulphonate	269-511-6	68259-08-5		
potassium perfluorohexane-1-sulphonate	223-393-2	3871-99-6		
perfluorohexane-1-sulphonic acid	206-587-1	355-46-4		
Ethanaminium, N-[4-[[4-(diethylamino)phenyl][4-(ethylamino)-1-naphthalenyl]methylene]-2,5-cyclohexadien-1-ylidene]-N-ethyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	1310480-24-0		
Methanaminium, N-[4-[[4-(dimethylamino)phenyl][4-(ethylamino)-1-naphthalenyl]methylene]-2,5-cyclohexadien-1-ylidene]-N-methyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	1310480-27-3		
Methanaminium, N-[4-[[4-(dimethylamino)phenyl][4-(phenylamino)-1-naphthalenyl]methylene]-2,5-cyclohexadien-1-ylidene]-N-methyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	1310480-28-4		
Beta-Cyclodextrin, compd. with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid ion(1-)(1:1)	-	1329995-45-0		
Gamma-Cyclodextrin, compd. with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid ion(1-)(1:1)	-	1329995-69-8		
Sulfonium, (thiodi-4,1-phenylene)bis[diphenyl-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid (1:2)	-	421555-73-9		
Iodonium, bis[4-(1,1-dimethylpropyl)phenyl]-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic	-	421555-74-0		
Sulfonium, tris[4-(1,1-dimethylethyl)phenyl]-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	425670-70-8		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, lithium salt (1:1)	-	55120-77-9		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, zinc salt	-	70136-72-0		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, compd. with N,N-diethylethanamine (1:1)	-	72033-41-1		
Iodonium, bis[(1,1-dimethylethyl)phenyl]-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid (1:1) (9Cl)	-	866621-50-3		
Sulfonium, (4-methylphenyl)diphenyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	910606-39-2		
Sulfonium, [4-[(2-methyl-1-oxo-2-propen-1-yl)oxy]phenyl]diphenyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	911027-68-4		

Sulfonium, [4-[(2-methyl-1-oxo-2-propenyl)oxy]phenyl]diphenyl-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid (1:1), polymer with 2-ethyltricyclo[3.3.1.13,7]dec-2-yl 2-methyl-2-propenoate, 3-hydroxytricyclo[3.3.1.13,7]dec-1-yl 2-methyl-2-propenoate and tetrahydro-2-oxo-3-furanyl 2-methyl-2-propenoate	-	911027-69-5		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, cesium salt (1:1)	-	92011-17-1		
Dibenzo[k,n][1,4,7,10,13]tetraoxathiacyclopentadecinium, 19-[4-(1,1-dimethylethyl)phenyl]-6,7,9,10,12,13-hexahydro-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	928049-42-7		
Sulfonium, triphenyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	144116-10-9		
Quinolinium, 1-(carboxymethyl)-4-[2-[4-[4-(2,2-diphenylethenyl)phenyl]-1,2,3,3a,4,8b-hexahydrocyclopent[b]indol-7-yl]ethenyl]-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	1462414-59-0		
Iodonium, diphenyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	153443-35-7		
Methanaminium, N,N,N-trimethyl-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid (1:1)	-	189274-31-5		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, compd.with 2-methyl-2-propanamine (1:1)	-	202189-84-2		
Iodonium, bis[4-(1,1-dimethylethyl)phenyl]-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	213740-81-9		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, gallium salt (9Cl)	-	341035-71-0		
Sulfonium, bis(4-methylphenyl)phenyl-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	341548-85-4		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, scandium(3+) salt (3:1)	-	350836-93-0		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, neodymium(3+) salt (3:1)	-	41184-65-0		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, yttrium(3+) salt (3:1)	-	41242-12-0		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, compd. With pyrrolidine (1:1)	-	1187817-57-7		
Ethanaminium, N,N,N-triethyl-, salt with1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid (1:1)	-	108427-55-0		
1-Butanaminium, N,N,N-tributyl-, salt with 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonic acid	-	108427-54-9		
Phosphonium, triphenyl(phenylmethyl)-, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-1-hexanesulfonate (1:1)	-	1000597-52-3		
1-Hexanesulfonic acid, 1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluoro-, sodium salt	-	82382-12-5		
p-(1,1-dimethylpropyl)phenol			201-280-9	80-46-6
Nonadecafluorodecanoic acid (PFDA) and its sodium and ammonium salts			-	-
Nonadecafluorodecanoic acid	206-400-3	335-76-2		
sodium nonadecafluorodecanoate	-	3830-45-3		
Decanoic acid, nonadecafluoro-, sodium salt	-	3830-45-3		
Ammonium nonadecafluorodecanoate	221-470-5	3108-42-7		

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			-	-
	4-heptylphenol	217-862-0	1987-50-4	
	Phenol, heptyl derivs.	276-743-1	72624-02-3	
	Phenol, 4-(1,1-diethylpropyl)-	-	37872-24-5	
	Phenol, 4-(1,1-dimethylpentyl)-	-	30784-31-7	
	Phenol, 4-(1,2,2-trimethylbutyl)-	-	911371-06-7	
	Phenol, 4-(1,2-dimethylpentyl)-	-	854904-93-1	
	Phenol, 4-(1,3,3-trimethylbutyl)-	-	911371-07-8	
	Phenol, 4-(1,3-dimethylpentyl)-	-	71945-81-8	
	Phenol, 4-(1,4-dimethylpentyl)-	-	857629-71-1	
	Phenol, 4-(1-ethyl-2,2-dimethylpropyl)-	-	861010-65-3	
	Phenol, 4-(1-ethyl-1,2-dimethylpropyl)-	-	30784-27-1	
	Phenol, 4-(1-ethylpentyl)-	-	6465-74-3	
	Phenol, 4-(1-methylhexyl)-	-	6863-24-7	
	Phenol, 4-(1-propylbutyl)-	-	6465-71-0	
	Phenol, 4-(3-ethylpentyl)-	-	911370-98-4	
	Phenol, 4-(3-methylhexyl)-	-	102570-52-5	
	Phenol, 4-(4-methylhexyl)-	-	1139800-98-8	
	Phenol, 4-(5-methylhexyl)-	-	100532-36-3	
	Phenol, 4-[2-methyl-1-(1-methylethyl)propyl]-	-	1824346-00-0	
	4-(2,3-dimethylpentan-2-yl)phenol	854-135-2	861011-60-1	
	4-(3-methylhexan-3-yl)phenol	854-958-7	30784-32-8	
	Phenol, 4-tert-heptyl-	-	288864-02-8	
	Phenol, 4-(1,1,3-trimethylbutyl)-	-	33104-11-9	
	Phenol, 4-(1,1,2,2-tetramethylpropyl)-	-	72861-06-4	
	Phenol, 4-(1-ethyl-3-methylbutyl)-	-	854904-92-0	
4,4'-isopropylidenediphenol Bisphenol A; BPA			201-245-8	80-05-7
Benzo[def]chrysene (Benzo[a]pyrene)			200-028-5	50-32-8
Perfluorononan-1-oic-acid and its sodium and ammonium salts			-	-
	Perfluorononan-1-oic-acid	206-801-3	375-95-1	
	Sodium salts of perfluorononan-1-oic-acid	-	-, 21049-39-8	
	Ammonium salts of perfluorononan-1-oic-acid	-	-, 4149-60-4	
Nitrobenzene			202-716-0	98-95-3
2-(2H-benzotriazol-2-yl)-4-(tert-butyl)-6-(sec-butyl)phenol (UV-350)			253-037-1	36437-37-3
2,4-di-tert-butyl-6-(5-chlorobenzotriazol-2-yl)phenol (UV-327)			223-383-8	3864-99-1
1,3-propanesultone			214-317-9	1120-71-4
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			-	-
	Reaction mass of 5-sec-butyl-2-(2,4-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane and 5-sec-butyl-2-(4,6-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane	413-720-9	117933-89-8	
	1,3-Dioxane, 2-[(1S,2S)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, trans-	-	676367-06-9	
	5-sec-butyl-2-(2,4-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane	-	-	
	5-sec-butyl-2-(4,6-dimethylcyclohex-3-en-1-yl)-5-methyl-1,3-dioxane	-	-	
	1,3-Dioxane, 2-(2,4-dimethyl-3-cyclohexen-1-yl)-5-methyl-5-(1-methylpropyl)-	-	186309-28-4	
	1,3-Dioxane, 2-(2,4-dimethyl-3-cyclohexen-1-yl)-5-methyl-5-(1-methylpropyl)-	-	117933-89-8	
	1,3-Dioxane, 2-[(1R,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, cis-	-	676367-05-8	
	1,3-Dioxane, 2-[(1R,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, cis-rel-	-	343934-04-3	

1,3-Dioxane, 2-[(1R,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, trans-	-	676367-09-2		
1,3-Dioxane, 2-[(1R,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, trans-rel-	-	343934-05-4		
1,3-Dioxane, 2-[(1R,2S)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, cis-	-	676367-04-7		
1,3-Dioxane, 2-[(1R,2S)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, trans-	-	676367-08-1		
1,3-Dioxane, 2-[(1S,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, cis-	-	676367-03-6		
1,3-Dioxane, 2-[(1S,2R)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, trans-	-	676367-07-0		
1,3-Dioxane, 2-[(1S,2S)-2,4-dimethyl-3-cyclohexen-1-yl]-5-methyl-5-(1-methylpropyl)-, cis-	-	676367-02-5		
Reaction mass of 5-[(2R)-butan-2-yl]-2-[(1R,2R)-2,4-dimethylcyclohex-3-en-1-yl]-5-methyl-1,3-dioxane and 5-[(2R)-butan-2-yl]-2-[(1R,6R)-4,6-dimethylcyclohex-3-en-1-yl]-5-methyl-1,3-dioxane and 5-[(2S)-butan-2-yl]-2-[(1R,2R)-2,4-dimethylcyclohex-3-en-1-yl]-5-methyl-1,3-dioxane and 5-[(2S)-butan-2-yl]-2-[(1S,2R)-2,4-dimethylcyclohex-3-en-1-yl]-5-methyl-1,3-dioxane and 5-[(2S)-butan-2-yl]-2-[(1S,6R)-4,6-dimethylcyclohex-3-en-1-yl]-5-methyl-1,3-dioxane	-	-		
1,2-benzenedicarboxylic acid, di-C6-10-alkyl esters or mixed decyl and hexyl and octyl diesters with ≥ 0.3% of dihexyl phthalate (EC No. 201-559-5)			-	-
1,2-Benzenedicarboxylic acid, mixed decyl and hexyl and octyl diesters	272-013-1	68648-93-1		
1,2-Benzenedicarboxylic acid, di-C6-10-alkyl esters	271-094-0	68515-51-5		
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)			-	-
Cadmium sulphate			233-331-6	10124-36-4, 31119-53-6
Cadmium fluoride			232-222-0	7790-79-6
2-ethylhexyl 10-ethyl-4,4-dioctyl-7-oxo-8-oxa-3,5-dithia-4-stannatetradecanoate (DOTE)			239-622-4	15571-58-1
2-benzotriazol-2-yl-4,6-di-tert-butylphenol (UV-320)			223-346-6	3846-71-7
2-(2H-benzotriazol-2-yl)-4,6-ditertpentylphenol (UV-328)			247-384-8	25973-55-1
Sodium peroxometaborate			231-556-4	7632-04-4
Sodium perborate, perboric acid, sodium salt			-	-
Perboric acid (H3BO2(O2)), monosodium salt, trihydrate	603-902-8	13517-20-9		
Borate(2-), tetrahydroxybis[μ-(peroxy-κO1:κO2)]di-, sodium (1:2)	-	90568-23-3		
Sodium perborate	239-172-9	15120-21-5		
Perboric acid, sodium salt	234-390-0	11138-47-9		
Borate(2-), tetrahydroxybis[μ-(peroxy-κO1:κO2)]di-, sodium, hydrate (1:2:6)	-	125022-34-6		
Cadmium chloride			233-296-7	10108-64-2
1,2-Benzenedicarboxylic acid, dihexyl ester, branched and linear			271-093-5	68515-50-4
Trixylyl phosphate			246-677-8	25155-23-1
Lead di(acetate)			206-104-4	301-04-2
Imidazolidine-2-thione (2-imidazoline-2-thiol)			202-506-9	96-45-7
Disodium 4-amino-3-[[4'-[(2,4-diaminophenyl)azo][1,1'-biphenyl]-4-yl]azo] -5-hydroxy-6-(phenylazo) naphthalene-2,7-disulphonate (C.I. Direct Black 38)			217-710-3	1937-37-7
Disodium 3,3'-[[1,1'-biphenyl]-4,4'-diylbis(azo)]bis(4-aminonaphthalene-1-sulphonate) (C.I. Direct Red 28)			209-358-4	573-58-0
Dihexyl phthalate			201-559-5	84-75-3
Cadmium sulphide			215-147-8	1306-23-6
Pentadecafluorooctanoic acid (PFOA)			206-397-9	335-67-1
Dipentyl phthalate (DPP)			205-017-9	131-18-0
Cadmium oxide			215-146-2	1306-19-0
Cadmium			231-152-8	7440-43-9
Ammonium pentadecafluorooctanoate (APFO)			223-320-4	3825-26-1

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			-	-
	Nonylphenol, ethoxylated	500-024-6	9016-45-9	
	4-Nonylphenol, ethoxylated 1 - 2.5 moles ethoxylated	500-045-0	26027-38-3	
	2-[2-[2-[2-(4-nonylphenoxy)ethoxy]ethoxy]ethoxy] ethanol	230-770-5	7311-27-5	
	2-[2-(4-nonylphenoxy)ethoxy]ethanol	243-816-4	20427-84-3	
	20-(4-nonylphenoxy)-3,6,9,12,15,18-hexaoxaicosan-1-ol	248-743-1	27942-27-4	
	26-(4-Nonylphenoxy)-3,6,9,12,15,18,21,24- octa-oxahexacosan -1-ol	-	14409-72-4	
	Poly(oxy-1,2-ethanediyl), α-(nonylphenyl)-ω-hydroxy-, branched	-	68412-54-4	
	Poly (oxy-1,2-ethanediyl), alpha -(nonylphenyl)-omega-hydroxy-, branched (CAS# 68412-54-4)	932-337-2	68412-54-4	
	Nonylphenol, ethoxylated (15-EO) (9016-45-9)	931-756-8	-	
	Nonylphenol, ethoxylated (10-EO) (9016-45-9)	931-755-2	-	
	Nonylphenol, ethoxylated (8-EO) (9016-45-9)	931-754-7	-	
	Nonylphenol, ethoxylated (6,5-EO) (9016-45-9)	931-753-1	-	
	Nonylphenol, branched, ethoxylated 1 - 2.5 moles ethoxylated	500-209-1	68412-54-4	
	2-[4-(3,6-dimethylheptan-3-yl)phenoxy]ethanol	687-832-3	1119449-37-4	
	14-(nonylphenoxy)-3,6,9,12-tetraoxatetradecan-1-ol	247-555-7	26264-02-8	
	Nonylphenol, ethoxylated (polymer)	938-618-6	-	
	Nonylphenol, ethoxylated (EO = 4)	939-975-0	-	
	Nonylphenol, ethoxylated (EO = 10)	939-993-9	-	
	Nonylphenolpolyglycolether	932-998-7	-	
	26-(nonylphenoxy)-3,6,9,12,15,18,21,24-octaoxahexacosan-1-ol	247-816-5	26571-11-9	
	Ethanol, 2-(4-nonylphenoxy)-	-	104-35-8	
	4-Nonylphenol, branched, ethoxylated (CAS: 127087-87-0)	932-098-4	127087-87-0	
	3,6,9,12-Tetraoxatetradecan-1-ol, 14-(4-nonylphenoxy)-	-	20636-48-0	
	4-t-Nonylphenol-diethoxylate	-	156609-10-8	
	Isononylphenol, ethoxylated	609-346-2	37205-87-1	
	p-Nonylphenol hexaethoxylate	-	34166-38-6	
	Poly(oxy-1,2-ethanediyl), a-(nonylphenyl)-w-hydroxy- (CAS 9016-45-9)	931-562-3	9016-45-9	
	4-Nonylphenol, branched, ethoxylated 1 - 2.5 moles ethoxylated	500-315-8	127087-87-0	
	23-(nonylphenoxy)-3,6,9,12,15,18,21-heptaoxatrico-san-1-ol	248-293-6	27177-05-5	
	2-{2-[4-(3,6-dimethylheptan-3-yl)phenoxy]ethoxy} ethanol	687-833-9	1119449-38-5	
Trilead dioxide phosphonate			235-252-2	12141-20-7
Trilead bis(carbonate) dihydroxide			215-290-6	1319-46-6
Tricosafluorododecanoic acid			206-203-2	307-55-1
Tetralead trioxide sulphate			235-380-9	12202-17-4
Tetraethyllead			201-075-4	78-00-2
Sulfurous acid, lead salt, dibasic			263-467-1	62229-08-7
Silicic acid, lead salt			234-363-3	11120-22-2
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			272-271-5	68784-75-8
Pyrochlore, antimony lead yellow -			232-382-1	8012-00-8
Pentalead tetraoxide sulphate			235-067-7	12065-90-6

Pentacosafluorotridecanoic acid			276-745-2	72629-94-8
Orange lead (lead tetroxide)			215-235-6	1314-41-6
o-toluidine			202-429-0	95-53-4
o-aminoazotoluene			202-591-2	97-56-3
n-pentyl-isopentylphthalate			933-378-9	776297-69-9
N-methylacetamide			201-182-6	79-16-3
N,N-dimethylformamide			200-679-5	68-12-2
Methyloxirane (Propylene oxide)			200-879-2	75-56-9
Methoxyacetic acid			210-894-6	625-45-6
Lead titanium zirconium oxide			235-727-4	12626-81-2
Lead titanium trioxide			235-038-9	12060-00-3
Lead oxide sulfate			234-853-7	12036-76-9
Lead monoxide (lead oxide)			215-267-0	1317-36-8
Lead dinitrate			233-245-9	10099-74-8
Lead cyanamidate			244-073-9	20837-86-9
Lead bis(tetrafluoroborate)			237-486-0	13814-96-5
Hexahydromethylphthalic anhydride including cis- and trans- stereo isomeric forms and all possible combinations of the isomers			-	-
Hexahydromethylphthalic anhydride	247-094-1	25550-51-0		
Hexahydro-3-methylphthalic anhydride	260-566-1	57110-29-9		
Hexahydro-1-methylphthalic anhydride	256-356-4	48122-14-1		
Hexahydro-4-methylphthalic anhydride	243-072-0	19438-60-9		
Heptacosafluorotetradecanoic acid			206-803-4	376-06-7
Henicosafluoroundecanoic acid			218-165-4	2058-94-8
Furan			203-727-3	110-00-9
Fatty acids, C16-18, lead salts			292-966-7	91031-62-8
Dioxobis(stearato)trilead			235-702-8	12578-12-0
Dinoseb (6-sec-butyl-2,4-dinitrophenol)			201-861-7	88-85-7
Dimethyl sulphate			201-058-1	77-78-1
Diisopentyl phthalate			210-088-4	605-50-5
Diethyl sulphate			200-589-6	64-67-5
Dibutyltin dichloride (DBTC)			211-670-0	683-18-1
Diazene-1,2-dicarboxamide (C,C'-azodi(formamide)) (ADCA)			204-650-8	123-77-3
Cyclohexane-1,2-dicarboxylic anhydride all possible combinations of the cis- and trans-isomers			-	-
cis-cyclohexane-1,2-dicarboxylic anhydride	236-086-3	13149-00-3		
trans-cyclohexane-1,2-dicarboxylic anhydride	238-009-9	14166-21-3		
Cyclohexane-1,2-dicarboxylic anhydride	201-604-9	85-42-7		
Bis(pentabromophenyl) ether (decabromodiphenyl ether) (DecaBDE)			214-604-9	1163-19-5
Biphenyl-4-ylamine			202-177-1	92-67-1
Acetic acid, lead salt, basic			257-175-3	51404-69-4
[Phthalato(2-)]dioxotrilead			273-688-5	69011-06-9
6-methoxy-m-toluidine (p-cresidine)			204-419-1	120-71-8
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			-	-
p-(1-methyloctyl)phenol	241-427-4	17404-66-9		
p-nonylphenol	203-199-4	104-40-5		
4-(1-Ethyl-1,4-dimethylpentyl)phenol	-	142731-63-3		
4-(1-Ethyl-1,3-dimethylpentyl)phenol	-	186825-36-5		
p-(1,1-dimethylheptyl)phenol	250-339-5	30784-30-6		
4-(1-ethyl-1-methylhexyl)phenol	257-907-1	52427-13-1		
p-isononylphenol	247-770-6	26543-97-5		
Phenol, 4-nonyl-, branched	284-325-5	84852-15-3		
4-(1,1,5-Trimethylhexyl)phenol	-	521947-27-3		
Isononylphenol	234-284-4	11066-49-2		

Phenol, nonyl-, branched	291-844-0	90481-04-2		
4-(3-ethylheptan-2-yl)phenol	635-696-0	186825-39-8		
Nonylphenol	246-672-0	25154-52-3		
4-methyl-m-phenylenediamine (toluene-2,4-diamine)			202-453-1	95-80-7
4-aminoazobenzene			200-453-6	60-09-3
4-(1,1,3,3-tetramethylbutyl)phenol, ethoxylated covering well-defined substances and UVCB substances, polymers and homologues			-	-
Poly(oxy-1,2-ethanediyl), α-[(1,1,3,3-tetramethylbutyl) phenyl]-ω-hydroxy-	-	9036-19-5		
Polyethylene glycol p-(1,1,3,3-tetramethylbutyl) phenyl ether	-	9002-93-1		
20-[4-(1,1,3,3-tetramethylbutyl)phe-noxy]-3,6,9,12,15,18-hexaoxaicosan-1-ol	219-682-8	2497-59-8		
2-[4-(1,1,3,3-tetramethylbutyl)phenoxy]ethanol	-	2315-67-5		
2-[2-[4-(1,1,3,3-tetramethylbutyl)phenoxy]ethoxy] ethanol, 2-{2-[4-(2,4,4-trimethylpentan-2-yl)phenoxy] ethoxy}ethanol	621-341-7	2315-61-9		
4,4'-oxydianiline and its salts			-	-
4,4'-oxydianiline	202-977-0	101-80-4		
4,4'-methylenedi-o-toluidine			212-658-8	838-88-0
3-ethyl-2-methyl-2-(3-methylbutyl)-1,3-oxazolidine			421-150-7	143860-04-2
1-bromopropane (n-propyl bromide)			203-445-0	106-94-5
1,2-diethoxyethane			211-076-1	629-14-1
1,2-Benzenedicarboxylic acid, dipentyl ester, branched and linear			284-032-2	84777-06-0
α,α-Bis[4-(dimethylamino)phenyl]-4 (phenylamino)naphthalene-1-methanol (C.I. Solvent Blue 4) with ≥ 0.1% of Michler's ketone (EC No. 202-027-5) or Michler's base (EC No. 202-959-2)			229-851-8	6786-83-0
N,N,N',N'-tetramethyl-4,4'-methylenedianiline (Michler's base)			202-959-2	101-61-1
Lead(II) bis(methanesulfonate) -			401-750-5	17570-76-2
Formamide			200-842-0	75-12-7
Diboron trioxide			215-125-8	1303-86-2
[4-[[4-anilino-1-naphthyl][4-(dimethylamino)phenyl]methylene]cyclohexa-2,5-dien-1-ylidene] dimethylammonium chloride (C.I. Basic Blue 26) with ≥ 0.1% of Michler's ketone (EC No. 202-027-5) or Michler's base (EC No. 202-959-2)			219-943-6	2580-56-5
[4-[4,4'-bis(dimethylamino) benzhydrylidene]cyclohexa-2,5-dien-1-ylidene]dimethylammonium chloride (C.I. Basic Violet 3) with ≥ 0.1% of Michler's ketone (EC No. 202-027-5) or Michler's base (EC No. 202-959-2)			208-953-6	548-62-9
4,4'-bis(dimethylamino)benzophenone (Michler's ketone)			202-027-5	90-94-8
4,4'-bis(dimethylamino)-4''-(methylamino)trityl alcohol with ≥ 0.1% of Michler's ketone (EC No. 202-027-5) or Michler's base (EC No. 202-959-2)			209-218-2	561-41-1
1,3,5-tris[(2S and 2R)-2,3-epoxypropyl]-1,3,5-triazine-2,4,6-(1H,3H,5H)-trione (β-TGIC) -			423-400-0	59653-74-6
1,3,5-Tris(oxiran-2-ylmethyl)-1,3,5-triazinane-2,4,6-trione (TGIC)			219-514-3	2451-62-9
1,2-bis(2-methoxyethoxy)ethane (TEGDME; triglyme)			203-977-3	112-49-2
1, 2-dimethoxyethane; ethylene glycol dimethyl ether (EGDME)			203-794-9	110-71-4
Zirconia Aluminosilicate Refractory Ceramic Fibres are fibres covered by index number 650-017-00-8 in Annex VI, part 3, table 3.1 of Regulation (EC) No 1272/2008 of the European Parliament and of the Council of 16 December 2008 on classification, labelling and packaging of substances and mixtures, and fulfil the three following conditions: a) oxides of aluminium, silicon and zirconium are the main components present (in the fibres) within variable concentration ranges b) fibres have a length weighted geometric mean diameter less two standard geometric errors of 6 or less micrometres (μm). c) alkaline oxide and alkali earth oxide (Na2O+K2O+CaO+MgO+BaO) content less or equal to 18% by weight			-	-
Refractories, fibers, aluminosilicate	-	142844-00-6		
Trilead diarsenate			222-979-5	3687-31-8
Potassium hydroxyoctaoxodizincatedichromate			234-329-8	11103-86-9
Phenolphthalein			201-004-7	77-09-8
Pentazinc chromate octahydroxide			256-418-0	49663-84-5
N,N-dimethylacetamide			204-826-4	127-19-5
Lead styphnate			239-290-0	15245-44-0
Lead dipicrate			229-335-2	6477-64-1
Lead diazide, Lead azide			236-542-1	13424-46-9

Formaldehyde, oligomeric reaction products with aniline			500-036-1	25214-70-4
Dichromium tris(chromate)			246-356-2	24613-89-6
Calcium arsenate			231-904-5	7778-44-1
Bis(2-methoxyethyl) phthalate			204-212-6	117-82-8
Bis(2-methoxyethyl) ether			203-924-4	111-96-6
Arsenic acid			231-901-9	7778-39-4
Aluminosilicate Refractory Ceramic Fibres are fibres covered by index number 650-017-00-8 in Annex VI, part 3, table 3.1 of Regulation (EC) No 1272/2008 of the European Parliament and of the Council of 16 December 2008 on classification, labelling and packaging of substances and mixtures, and fulfil the three following conditions: a) oxides of aluminium and silicon are the main components present (in the fibres) within variable concentration ranges b) fibres have a length weighted geometric mean diameter less two standard geometric errors of 6 or less micrometres (µm) c) alkaline oxide and alkali earth oxide (Na2O+K2O+CaO+MgO+BaO) content less or equal to 18% by weight			-	-
	Refractories, fibers, aluminosilicate	-	142844-00-6	
4-(1,1,3,3-tetramethylbutyl)phenol			205-426-2	140-66-9
2-Methoxyaniline, o-Anisidine			201-963-1	90-04-0
2,2'-dichloro-4,4'-methylenedianiline			202-918-9	101-14-4
1,2-dichloroethane			203-458-1	107-06-2
Strontium chromate			232-142-6	7789-06-2
Hydrazine			206-114-9	302-01-2, 7803-57-8
2-ethoxyethyl acetate			203-839-2	111-15-9
1-Methyl-2-pyrrolidone (NMP)			212-828-1	872-50-4
1,2-Benzenedicarboxylic acid, di-C7-11-branched and linear alkyl esters			271-084-6	68515-42-4
1,2-Benzenedicarboxylic acid, di-C6-8-branched alkyl esters, C7-rich			276-158-1	71888-89-6
1,2,3-trichloropropane			202-486-1	96-18-4
Cobalt(II) sulphate			233-334-2	10124-43-3
Cobalt(II) dinitrate			233-402-1	10141-05-6
Cobalt(II) diacetate			200-755-8	71-48-7
Cobalt(II) carbonate			208-169-4	513-79-1
Chromium trioxide			215-607-8	1333-82-0
Acids generated from chromium trioxide and their oligomers			-	-
	Oligomers of chromic acid and dichromic acid	-	-	
	Chromic acid	231-801-5	7738-94-5	
	Dichromic acid	236-881-5	13530-68-2	
2-methoxyethanol			203-713-7	109-86-4
2-ethoxyethanol			203-804-1	110-80-5
Trichloroethylene			201-167-4	79-01-6
Tetraboron disodium heptaoxide, hydrate			235-541-3	12267-73-1
Sodium chromate			231-889-5	7775-11-3
Potassium dichromate			231-906-6	7778-50-9
Potassium chromate			232-140-5	7789-00-6
Disodium tetraborate, anhydrous			215-540-4	12179-04-3, 1303-96-4, 1330-43-4
Boric acid EC No. 233-139-2 and EC No. 234-343-4			-	-
	Boric acid, crude natural	234-343-4	11113-50-1	
	Boric acid	233-139-2	10043-35-3	
Ammonium dichromate			232-143-1	7789-09-5
Acrylamide			201-173-7	79-06-1
Tris(2-chloroethyl) phosphate			204-118-5	115-96-8
Pitch, coal tar, high-temp. -			266-028-2	65996-93-2
Lead sulfochromate yellow (C.I. Pigment Yellow 34) -			215-693-7	1344-37-2
Lead chromate molybdate sulphate red (C.I. Pigment Red 104) -			235-759-9	12656-85-8
Lead chromate			231-846-0	7758-97-6
Diisobutyl phthalate			201-553-2	84-69-5

Anthracene oil, anthracene-low -			292-604-8	90640-82-7
Anthracene oil, anthracene paste, distn. lights -			295-278-5	91995-17-4
Anthracene oil, anthracene paste, anthracene fraction -			295-275-9	91995-15-2
Anthracene oil, anthracene paste -			292-603-2	90640-81-6
Anthracene oil -			292-602-7	90640-80-5
2,4-dinitrotoluene			204-450-0	121-14-2
Triethyl arsenate -			427-700-2	15606-95-8
Sodium dichromate			234-190-3	10588-01-9, 7789-12-0
Lead hydrogen arsenate			232-064-2	7784-40-9
Hexabromocyclododecane (HBCDD) and all major diastereoisomers identified			-	-
	Hexabromocyclododecane	247-148-4	25637-99-4	
	1,2,5,6,9,10-hexabromocyclododecane	221-695-9	3194-55-6	
	alpha-hexabromocyclododecane	-	134237-50-6	
	beta-hexabromocyclododecane	-	134237-51-7	
	gamma-hexabromocyclododecane	-	134237-52-8	
Dibutyl phthalate (DBP)			201-557-4	84-74-2
Diarsenic trioxide			215-481-4	1327-53-3
Diarsenic pentaoxide			215-116-9	1303-28-2
Cobalt dichloride			231-589-4	7646-79-9
Bis(tributyltin) oxide (TBTO)			200-268-0	56-35-9
Bis (2-ethylhexyl)phthalate (DEHP)			204-211-0	117-81-7
Benzyl butyl phthalate (BBP)			201-622-7	85-68-7
Anthracene			204-371-1	120-12-7
Alkanes, C10-13, chloro (Short Chain Chlorinated Paraffins)			287-476-5	85535-84-8
5-tert-butyl-2,4,6-trinitro-m-xylene (Musk xylene)			201-329-4	81-15-2
4,4'- Diaminodiphenylmethane (MDA)			202-974-4	101-77-9

Extrait 17 - **Règlementation REACH**
Restrictions en matière de fabrication, de mise sur le marché et d'utilisation de certaines substances, mélanges et articles dangereux

https://echa.europa.eu/it/substances-restricted-under-reach?p_p_id=disslists_WAR_disslistsportlet&p_p_lifecycle=0&p_p_state=normal&p_p_mode=view&_disslists_WAR_disslistsportlet_cur=1&_disslists_WAR_disslistsportlet_substance_identifier_field_key=&_disslists_WAR_disslistsportlet_delta=50&_disslists_WAR_disslistsportlet_doSearch=&_disslists_WAR_disslistsportlet_prc_entry_no=&_disslists_WAR_disslistsportlet_deltaParamValue=50&_disslists_WAR_disslistsportlet_orderByCol=prc_entry_no&_disslists_WAR_disslistsportlet_orderByType=desc

FINDER déclare que dans ses propres produits, les substances objets de la liste de l'annexe XVII de la réglementation REACH, ne sont pas présentes en concentration supérieure aux limites imposées.