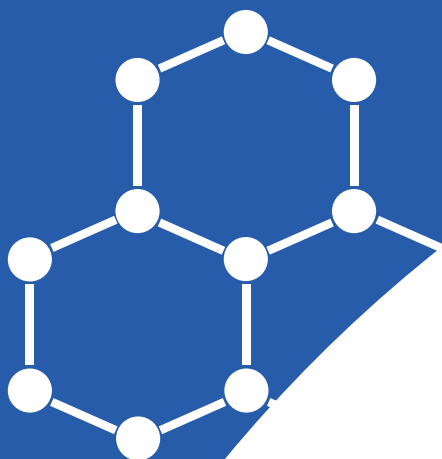




finder[®]
SWITCH TO THE FUTURE

REACH - Reglamento (UE) n° 1907/2006

REACH es un reglamento comunitario sobre las sustancias químicas, sobre su control y su empleo seguro. Corresponde al registro, a la valoración, a la autorización y a la restricción de las sustancias químicas (REACH es el acrónimo de **R**egistration, **E**valuation, **A**uthorisation and **R**estriction of **C**hemicals).

REACH sustituye numerosas normas comunitarias relativas a las sustancias químicas (también como componentes individuales presentes en preparados) y desarrolla una función complementaria con respecto a otras normativas ambientales y de seguridad.

Uno de los objetivos más importantes es, sin duda, la tutela de la salud humana y del medio ambiente respecto a los riesgos que se podrían originar por el empleo de sustancias químicas en cuanto tales o en cuanto componentes de un preparado o de un artículo.

Estas restricciones entran en un contexto que Finder está tomando en consideración desde hace tiempo. Como **usuario final (posterior)** de sustancias químicas, mantenemos activa la comunicación con nuestros proveedores, analizando y teniendo bajo estrecho control las fichas de seguridad previstas para cada tipo de sustancias o preparados. Se recuerda que en sus propios procesos productivos Finder emplea sustancias y/o preparados que no modifican las características químicas de los mismos.

En los distintos plazos previstos por la Directiva, Finder se compromete a prestar una atención especial a la selección de los proveedores, dando como garantía la comunicación con sus propios clientes, de manera más específica por lo que respecta a las actualizaciones de las tablas que divulgan los organismos oficiales (ECHA) - A continuación sigue la lista de las sustancias consideradas altamente peligrosas "Candidate list of substances of very high concern for authorization (SVHC)" y la extrapolación de la lista de sustancias del Anexo XVII del Reglamento Reach actualizado.

Además, la política de Finder ha sido implementada siguiendo el sistema ISO 14001 (Sistema de Gestión Ambiental) que promueve cuestionarios destinados a evaluar, con una finalidad preventiva, los sujetos con los que colabora. Siendo **usuarios finales** de sustancias químicas, se aplican medidas de precaución de los riesgos que se podrían derivar de las mismas. Por este motivo, además de lo ya mencionado, se ha puesto en marcha un Sistema de Gestión de Salud y Seguridad que garantiza la tutela de las personas implicadas en los distintos procesos de producción.

REACH – Sustancias altamente preocupantes
Candidate List of Substances of Very High Concern – SVHC
<https://echa.europa.eu/candidate-list-table>

FINDER informa que en sus productos no se encuentran las sustancias que se indican en la tabla, en concentraciones superiores al 0,1% en peso.

Nombre de la sustancia	Número EC		Número EC	Número CAS
	Número EC	Número CAS		
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-diylbis(oxy)]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-butanefulfonate	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
1,6,7,8,9,14,15,16,17,18-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			-	-
1,6,7,8,9,14,15,16,17,18-dodecachloropentacyclo[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)-1,6,7,8,9,14,15,16,17,18,18-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) (9CI)	-	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[3.3.1.1 ^{3,7}]dec-2-yl 2-methyl-2-propenoate, 3-hydroxytricyclo[3.3.1.1 ^{3,7}]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-(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 with 1,1,2,2,3,3,4,4,5,5,6,6,6-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-(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), α -(nonylphenyl)- ω -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	-		
Nonylphenolpolyglycoether	932-998-7	-		
26-(nonylphenoxy)-3,6,9,12,15,18,21,24-octa-oxahexacosan-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), α -(nonylphenyl)- ω -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-hepta-oxatricosan-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		
Tricosafuorododecanoic 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 (H ₂ SiO ₅), 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		

Pentacosafuorotridecanoic 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
Heptacosafuorotetradecanoic acid	206-803-4	376-06-7
Henicosafuoroundecanoic 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)phenoxy]-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-triazine-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 (Na ₂ O+K ₂ O+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 (Na ₂ O+K ₂ O+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

Extracto del anexo 17 - Reglamento Reach

Restricciones en materia de fabricación, introducción en el mercado y empleo de sustancias, mezclas y artículos peligrosos

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 declara que en sus productos no están presentes las sustancias enumeradas en el anexo XVII del Reglamento Reach en concentraciones superiores a los límites prescritos.