The use of the following substances in the materials for the products manufactured and sold by the Sysmex group are prohibited or restricted.

|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 1 | White phosphorus | - | ISHA |
| 2 | Benzidine and its salt | 10000 | ISHA |
| 3 | 4-Aminodiphenyl and its salts | 10000 | ISHA |
| 4 | Asbestos | 1000 | ISHA |
| 5 | 4-nitrodiphenyl and its salts | 10000 | ISHA |
| 6 | Bis (chloromethyl) ether | 10000 | ISHA |
| 7 | $\beta$-Naphthylamine and its salt | 10000 | ISHA |
| 8 | A rubber paste containing benzene, the volume of benzene contained being more than $5 \%$ of the solvent (including the filler) of the rubber paste | - | ISHA |
| 9 | Polychlorinated biphenyl | - | ECS, POPs Convention |
| 10 | Naphthalene, chloro derivs. | - | ECS, POPs Convention |
| 11 | HEXACHLOROBENZENE | - | ECS, POPs Convention |
| 12 | ALDRIN | - | ECS, POPs Convention |
| 13 | DIELDRIN | - | ECS, POPs Convention |
| 14 | ENDRIN | - | ECS, POPs Convention |
| 15 | 4,4'-DDT | - | ECS, POPs Convention |
| 16 | CHLORDANE | - | ECS, POPs Convention |
| 17 | Bis(tributyltin) oxide (TBTO) | - | ECS, POPs Convention |
| 18 | DTPD OEKANAL ( N N'-DITOLYLPARAPHENYLENE | - | ECS, POPs Convention |
| 19 | 2,4,6-Tri-tert-butylphenol | - | ECS, POPs Convention |
| 20 | TOXAPHENE | - | ECS, POPs Convention |
| 21 | MIREX | - | ECS, POPs Convention |
| 22 | Dicofol | - | ECS, POPs Convention |
| 23 | Hexachloro-1,3-butadiene | - | ECS, POPs Convention |
| 24 | 2-(2H-1,2,3-Benzotriazol-2-yl)-4,6-di-tert-butylphenol | - | ECS, POPs Convention |
| 25 | HEPTADECAFLUOROOCTANESULFONIC ACID | - | ECS, POPs Convention |
| 26 | Perfluoro-1-octanesulfonyl fluoride | - | ECS, POPs Convention |
| 27 | PENTACHLOROBENZENE | - | ECS, POPs Convention |
| 28 | ALPHA-HCH | - | ECS, POPs Convention |
| 29 | BETA-HCH | - | ECS, POPs Convention |
| 30 | LINDANE | - | ECS, POPs Convention |
| 31 | KEPONE (TM) | - | ECS, POPs Convention |
| 32 | HEXABROMOBIPHENYL | - | ECS, POPs Convention |
| 33 | diphenyl ether, tetrabromo derivative | - | ECS, POPs Convention |
| 34 | Pentabromodiphenyl ether | - | ECS, POPs Convention |
| 35 | 2,2',4,4,5,5'-HEXABROMODIPHENYL ETHER 2,2,4,4',5,6'-HEXABROMODIPHENYL ETHER | - | ECS, POPs Convention |
| 36 | $\begin{aligned} & \text { PBDE } 175 \\ & \text { 2,2',3,4,4,'5,6-Heptabromodiphenyl Ether } \end{aligned}$ | - | ECS, POPs Convention |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 37 | Thiosulfan Endosulfan I BETA-ENDOSULFAN | - | ECS, POPs Convention |
| 38 | Hexabromocyclododecane | - | ECS, POPs Convention |
| 39 | Pentachlorophenol | - | ECS, POPs Convention |
| 40 | Polychlorinated normal paraffin (It is limited that the number of carbon is 10 to 13 and the content of chlorine is more than $48 \%$ of the total weight.) | - | ECS, POPs Convention |
| 41 | Decabromodiphenyl oxide | - | ECS, POPs Convention |
| 42 | Pentadecafluoro-n-octanoic acid | - | ECS, POPs Convention |
| 43 | Radioactive substances | - | The Nuclear Reactor Regulation Law |
| 44 | Shortchain Chlorinated Paraffins (C10-C13) | No intentional addition | POPs Convention |
| 45 | Polychlorinated dibenzo-p-dioxins (PCDD) | No intentional addition | POPs Convention |
| 46 | Polychlorinated dibenzofurans (PCDF) | No intentional addition | POPs Convention |
| 47 | trichlorofluoromethane(CFC11) | - | Montreal Protocol |
| 48 | Dichlorodifluoromethane(CFC12) | - | Montreal Protocol |
| 49 | Trichlorotrifluoroethane (CFC113) | - | Montreal Protocol |
| 50 | Dichlorotetrafluoroethane(CFC 114) | - | Montreal Protocol |
| 51 | chloropentafluoroethane(CFC 115) | - | Montreal Protocol |
| 52 | Bromochlorodifluoromethane(Halon 1211) | - | Montreal Protocol |
| 53 | Bromotrifluoromethane (Halon1301) | - | Montreal Protocol |
| 54 | Dibromotetrafluoroethane(Halon 2402) | - | Montreal Protocol |
| 55 | Chlorotrifluoromethane(CFC13) | - | Montreal Protocol |
| 56 | Pentachlorofluoroethane(CFC 111) | - | Montreal Protocol |
| 57 | Tetrachlorodifluoroethane(CFC 112) | - | Montreal Protocol |
| 58 | Heptachlorofluoropropane(CFC 211) | - | Montreal Protocol |
| 59 | Hexachlorodifluoropropane(CFC 212) | - | Montreal Protocol |
| 60 | Pentachlorotrifluoropropane(CFC 213) | - | Montreal Protocol |
| 61 | Tetrachlorotetrafluoropropane(CFC 214) | - | Montreal Protocol |
| 62 | Trichloropentafluoropropane(CFC 215) | - | Montreal Protocol |
| 63 | Dichlorohexafluoropropane(CFC 216) | - | Montreal Protocol |
| 64 | chloroheptafluoropropane(CFC 217) | - | Montreal Protocol |
| 65 | Carbon tetrachloride | - | Montreal Protocol |
| 66 | 1,1,1-trichloroethane | - | Montreal Protocol |
| 67 | Dichlorofluoromethane(HCFC 21) | - | Montreal Protocol |
| 68 | Chlorodifluoromethane(HCFC 22) | - | Montreal Protocol |
| 69 | Chlorofluoromethane (HCFC31) | - | Montreal Protocol |
| 70 | Tetrachlorofluoroethane(HCFC 121) | - | Montreal Protocol |
| 71 | Trichlorodifluoroethane(HCFC 122) | - | Montreal Protocol |
| 72 | Dichlorotrifluoroethane(HCFC 123) | - | Montreal Protocol |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 73 | 2,2-dichloro-1,1,1-trifluoroethane(HCFC 123) | - | Montreal Protocol |
| 74 | Chlorotetrafluoroethane (HCFC 124) | - | Montreal Protocol |
| 75 | 2-Chloro-1,1,1,2-tetrafluoroethane(HCFC 124) | - | Montreal Protocol |
| 76 | Trichlorofluoroethane (HCFC131) | - | Montreal Protocol |
| 77 | Dichlorodifluoroethane(HCFC 132) | - | Montreal Protocol |
| 78 | Chlorotrifluoroethane (HCFC133) | - | Montreal Protocol |
| 79 | Dichlorofluoroethane(HCFC141) | - | Montreal Protocol |
| 80 | 1,1-Dichloro-1-fluoroethane(HCFC-141b) | - | Montreal Protocol |
| 81 | Chlorodifluoroethane (HCFC142) | - | Montreal Protocol |
| 82 | 1-Chloro-1,1-difluoroethane(HCFC142b) | - | Montreal Protocol |
| 83 | Chlorofluoroethane(HCFC-151) | - | Montreal Protocol |
| 84 | Hexachlorofluoropropane(HCFC 221) | - | Montreal Protocol |
| 85 | Pentachlorodifluoropropane(HCFC 222) | - | Montreal Protocol |
| 86 | Tetrachlorotrifluoropropane(HCFC 223) | - | Montreal Protocol |
| 87 | Trichlorotetrafluoropropane(HCFC 224) | - | Montreal Protocol |
| 88 | Dichloropentafluoropropane(HCFC 225) | - | Montreal Protocol |
| 89 | 3,3-Dichloro-1,1,1,2,2-pentafluoropropane (HCFC225ca) | - | Montreal Protocol |
| 90 | 1,3-Dichloro-1,1,2,2,3-pentafluoropropane (HCFC225cb) | - | Montreal Protocol |
| 91 | Chlorohexafluoropropane(HCFC 226) | - | Montreal Protocol |
| 92 | Pentachlorofluoropropane(HCFC 231) | - | Montreal Protocol |
| 93 | Tetrachlorodifluoropropane(HCFC 232) | - | Montreal Protocol |
| 94 | Trichlorotrifluoropropane(HCFC 233) | - | Montreal Protocol |
| 95 | Dichlorotetrafluoropropane(HCFC 234) | - | Montreal Protocol |
| 96 | Chloropentafluoropropane(HCFC 235) | - | Montreal Protocol |
| 97 | Tetrachlorofluoropropane(HCFC 241) | - | Montreal Protocol |
| 98 | Trichlorodifluoropropane(HCFC 242) | - | Montreal Protocol |
| 99 | Dichlorotrifluoropropane(HCFC 243) | - | Montreal Protocol |
| 100 | Chlorotetrafluoropropane(HCFC 244) | - | Montreal Protocol |
| 101 | Trichlorofluoropropane(HCFC 251) | - | Montreal Protocol |
| 102 | Dichlorodifluoropropane(HCFC 252) | - | Montreal Protocol |
| 103 | Chlorotrifluoropropane(HCFC 253) | - | Montreal Protocol |
| 104 | Dichlorofluoropropane (HCFC261) | - | Montreal Protocol |
| 105 | Chlorodifluoropropane (HCFC262) | - | Montreal Protocol |
| 106 | Chlorofluoropropane (HCFC271) | - | Montreal Protocol |
| 107 | Dibromofluoromethane | - | Montreal Protocol |
| 108 | Bromodifluoromethane(HBFC 22B1) | - | Montreal Protocol |
| 109 | Bromofluoromethane | - | Montreal Protocol |
| 110 | Tetrabromofluoroethane | - | Montreal Protocol |
| 111 | Tribromodifluoroethane | - | Montreal Protocol |
| 112 | Dibromotrifluoroethane | - | Montreal Protocol |
| 113 | Bromo tetrafluoroethane | - | Montreal Protocol |
| 114 | Tribromofluoroethane | - | Montreal Protocol |
| 115 | Dibromodifluoroethane | - | Montreal Protocol |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 116 | Bromotrifluoroethane | - | Montreal Protocol |
| 117 | Dibromofluoroethane | - | Montreal Protocol |
| 118 | Bromodifluoroethane | - | Montreal Protocol |
| 119 | Bromofluoroethane (HBFC-151 B1) | - | Montreal Protocol |
| 120 | Hexabromofluoropropane | - | Montreal Protocol |
| 121 | Pentabromodifluoropropane | - | Montreal Protocol |
| 122 | Tetrabromotrifluoropropane | - | Montreal Protocol |
| 123 | Tribromotetrafluoropropane | - | Montreal Protocol |
| 124 | Dibromopentafluoropropane | - | Montreal Protocol |
| 125 | Bromohexafluoropropane | - | Montreal Protocol |
| 126 | Pentabromofluoropropane | - | Montreal Protocol |
| 127 | Tetrabromodifluoropropane | - | Montreal Protocol |
| 128 | Tribromotrifluoropropane | - | Montreal Protocol |
| 129 | Dibromotrifluoropropane | - | Montreal Protocol |
| 130 | Bromopentafluoropropane | - | Montreal Protocol |
| 131 | Tetrabromofluoropropane | - | Montreal Protocol |
| 132 | Tribromodifluoropropane | - | Montreal Protocol |
| 133 | Dibromotrifluoropropane | - | Montreal Protocol |
| 134 | Bromo tetrafluoropropane | - | Montreal Protocol |
| 135 | Tribromofluoropropane | - | Montreal Protocol |
| 136 | Dibromodifluoropropane | - | Montreal Protocol |
| 137 | Bromotrifluoropropane | - | Montreal Protocol |
| 138 | Dibromofluoropropane | - | Montreal Protocol |
| 139 | Bromodifluoropropane | - | Montreal Protocol |
| 140 | Bromofluoropropane | - | Montreal Protocol |
| 141 | Bromochloromethane | - | Montreal Protocol |
| 142 | Methylbromide | - | Montreal Protocol |
| 143 | O-Alkyl (<=C10 including cycloalkyl) alkyl (Me,Et,n-Pr or i-Pr)phosphonofluoridates | - | PCW |
| 144 | O-Alkyl (<=C10 including cycloalkyl) N,N-dialkyl (Me,Et,n-Pr or i-Pr) phosphoramidocyanidates | - | PCW |
| 145 | O-Alkyl (<=C10 including cycloalkyl) S-2-dialkyl (Me,Et,n-Pr or i-Pr)aminoethyl alkyl (Me,Et,n-Pr or i-Pr) phosphonothiolates and corresponding alkylated or protonated salts | - | PCW |
| 146 | S-2-dialkyl (Me,Et,n-Pr or i-Pr) aminoethyl hydrogen alkyl (Me,Et,n-Pr or i-Pr) phosphonothiolates and corresponding alkylated or protonated salts | - | PCW |
| 147 | 2-Chloroethylchloromethylsulfide | - | PCW |
| 148 | Bis(2-chloroethyl)sulfide (Mustard gas) | - | PCW |
| 149 | Bis(2-chloroethylthio)methane | - | PCW |
| 150 | 1,2-Bis(2-chloroethylthio)ethane (Sesquimustard) | - | PCW |
| 151 | 1,3-Bis(2-chloroethylthio)-n-propane | - | PCW |
| 152 | 1,4-Bis(2-chloroethylthio)-n-butane | - | PCW |
| 153 | 1,5-Bis(2-chloroethylthio)-n-pentane | - | PCW |
| 154 | Bis(2-chloroethylthiomethyl)ether | - | PCW |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 155 | Bis(2-chloroethylthioethyl)ether (O-Mustard) | - | PCW |
| 156 | 2-Chlorovinyldichloroarsine (Lewisite 1) | - | PCW |
| 157 | Bis(2-chlorovinyl)chloroarsine (Lewisite 2) | - | PCW |
| 158 | Tris(2-chlorovinyl)arsine (Lewisite 3) | - | PCW |
| 159 | Bis(2-chloroethyl)ethylamine (HN1) | - | PCW |
| 160 | Bis(2-chloroethyl)methylamine (HN2) | - | PCW |
| 161 | Tris(2-chloroethyl)amine (HN3) | - | PCW |
| 162 | Saxitoxin | - | PCW |
| 163 | Ricin | - | PCW |
| 164 | P-alkyl-N-[1-(dialkylamino)alkylidene]phosphonamidic acid=fluoride (P -Alkyl or dialkylamino alkyl group is a cycloalkyl group and the alkylidene group is a cycloalkylmethylidene group, and the number of carbons in the alkyl group of the P-alkyl or dialkylamino alkyl group is 10 or less and the number of carbons of the alkylidene group is 11 (where the alkylidene group is a cycloalkylmethylidene group, the number of carbons of the cycloalkyl alkyl group is 10) or less). and alkylated salts and protonated salts thereof. | - | PCW |
| 165 | N -[1-(dialkylamino)alkylidene] phosphonamidic acid=fluoride (including those in which the alkyl group is a cycloalkyl group and the alkylidene group is a cycloalkyl methylidene group, where the alkyl group (limited to those having not more than ten carbons and not more than eleven carbons of the alkylidene group (where the alkylidene group is a cycloalkylmethylidene group, the number of carbons of the alkyl group of the cycloalkyl is ten)). and alkylated salts and protonated salts thereof. | - | PCW |
| 166 | N -[1-(dialkylamino)alkylidene]phosphoramidofluoridic acid (including those in which the alkyl group is a cycloalkyl group and the alkylidene group is a cycloalkylmethylidene group, where the alkyl group (limited to those having not more than ten carbons and not more than eleven carbons of the alkylidene group (where the alkylidene group is a cycloalkyl methylidene group, the number of carbons of the alkyl group of the cycloalkyl group is ten)). and alkylated salts and protonated salts thereof. | - | PCW |
| 167 | Alkyl=N-[1-(dialkylamino)alkylidene]phosphoramidofluoridate (wherein the alkyl group attached to the phosphoramidofluoridate or the alkyl group of the dialkylamino is a cycloalkyl group) and those in which the alkylidene group is a cycloalkylmethylidene group, wherein the number of carbons of the alkyl group attached to the phosphoramido fluoride and the alkyl group of the dialkylamino is less than ten and the number of carbons of the alkylidene group is eleven (the alkylidene group is a cycloalkylmethylidene group) In the case where the alkyl group is a cycloalkyl methylidene group, the number of carbons of the alkyl group of the cycloalkyl group is not more than 10)). and alkylated salts and protonated salts thereof. | - | PCW |
| 168 | N -[bis(diethylamino)methylidene]-P-methylphosphonamide acid = Fluoride | - | PCW |
| 169 | N -Acetyloxyalkyl-N,N, $\mathrm{N}^{\prime}$, $\mathrm{N}^{\prime}$-tetraalkyl- $\mathrm{N}^{\prime}$-\{[3-(dimethylcarbamoyloxy)pyridin-2-yl]methyl\} $\mathrm{N}, \mathrm{N}^{\prime}$-(decan-1,Xdiyl)diammonium=dibromide (where the acetyloxyalkyl group is a cyanoalkyl group or Including one in which the acetyloxyalkyl (where the acetyloxyalkyl group is a cyanoalkyl or hydroxyalkyl group, respectively cyanoalkyl or hydroxyalkyl) and tetraalkyl alkyl groups have ten or less carbons, and (iii) an acetyloxy group (if the acetyloxyalkyl group is a cyanoalkyl or hydroxyalkyl group, a cyano group or a hydroxy | - | PCW |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
|  | group, respectively) is bonded to the alkyl group to which it is attached at any carbon atom in position number 1 to 8 of the alkyl group concerned. ( $X$ shall represent an integer number from 1 to 10.) |  |  |
| 170 | $\mathrm{N}, \mathrm{N}, \mathrm{N}^{\prime}, \mathrm{N}^{\prime}$-tetraalkyl-N, $\mathrm{N}^{\prime}$ ' -bis\{[3-(dimethylcarbamoyloxy)pyridin-2- <br> yl]methyl\}ethanbis(amidium)=dibromide (limited to those having 10 or less carbons in the tetraalkyl alkyl group) (limited to those with 10 or less alkyl carbons in the tetraalkyl group). | - | PCW |
| 171 | N,N,N', N' -tetraalkyl-N,N"'bis\{[3-(dimethylcarbamoyloxy)pyridine-2-yl]methyl\}-N,N, $\mathrm{N}^{\prime}$-(2,X1-dioxoalkane-1,X2-diyl)diammonium=. Dibromide (provided that the structure of the alkane is linear, the number of carbons of the alkane concerned is between four and twelve and the number of carbons of the alkyl group of the tetraalkyl is less than ten). (X1 shall represent the number of carbons of the alkane concerned minus one and $X 2$ shall represent a number equal to the number of carbons of the alkane concerned). | - | PCW |
| 172 | Alkyl ( $\mathrm{Me}, \mathrm{Et}, \mathrm{n}-\mathrm{Pr}$ or i-Pr) phosphonyldifluorides | - | PCW |
| 173 | O-Alkyl (<=C10 including cycloalkyl) O-2-dialkyl (Me,Et,n-Pr or i-Pr)aminoethyl alkyl (Me,Et,n-Pr or i-Pr) phosphonites and corresponding alkylated or protonated salts | - | PCW |
| 174 | O-2-Dialkyl (Me,Et,n-Pr or i-Pr) aminoethyl hydrogen alkyl (Me,Et,n-Pr or i-Pr) phosphonites and corresponding alkylated or protonated salts | - | PCW |
| 175 | O-Isopropyl methylphosphonochloridate (Chlorosarin) | - | PCW |
| 176 | O-Pinacolyl methylphosphonochloridate (Chlorosoman) | - | PCW |
| 177 | O,O-Diethyl S-[2-(diethylamino)ethyl] phosphorothiolate (also known as amiton) and corresponding alkylated or protonated salts | - | PCW |
| 178 | 1,1,3,3,3-Pentafluoro-2-(trifluoromethyl)-1-propene (PFIB) | - | PCW |
| 179 | 3-Quinuclidinyl benzilate (BZ) | - | PCW |
| 180 | Chemicals, except for those in the following, containing a phosphorus atom to which is bonded alkyl( $\mathrm{Me}, \mathrm{Et}, \mathrm{n}-\mathrm{Pr}$ or $\mathrm{i}-\mathrm{Pr}$ ) but is not bonded any other carbon atoms (i) Chemicals in Row 1, Column 3(1) to (4) and Column 4 (ii) O-Ethyl S-phenyl thylphosphonothiolothionate(Fonofos) | - | PCW |
| 181 | N,N-Dialkyl (Me,Et,n-Pr or i-Pr) phosphoramidic dihalides | - | PCW |
| 182 | Dialkyl (Me,Et,n-Pr or i-Pr) N,N-dialkyl (Me,Et,n-Pr or i-Pr) phosphoramidate | - | PCW |
| 183 | Arsenic trichloride | - | PCW |
| 184 | 2,2-Diphenyl-2-hydroxyacetic acid | - | PCW |
| 185 | Quinuclidin-3-ol | - | PCW |
| 186 | $\mathrm{N}, \mathrm{N}$-Dialkyl (Me,Et,n-Pr or i-Pr)aminoethyl-2-chlorides and corresponding protonated salts | - | PCW |
| 187 | N,N-Dialkyl (Me,Et,n-Pr or i-Pr)aminoethane-2-ols (excluding N,NDimethylaminoethanol, $\mathrm{N}, \mathrm{N}$-Diethylaminoethanol and corresponding protonated salts) | - | PCW |
| 188 | N,N-Dialkyl (Me,Et,n-Pr or i-Pr)aminoethane-2-thiols and corresponding protonated salts | - | PCW |
| 189 | Bis(2-hydroxyethyl)sulfide (Thiodiglycol) | - | PCW |
| 190 | 3,3-Dimethylbutan-2-ol (Pinacolyl alcohol) | - | PCW |
| 191 | Carbonyl dichloride (Phosgene) | - | PCW |
| 192 | Cyanogen chloride | - | PCW |
| 193 | Hydrogen cyanide | - | PCW |
| 194 | Trichloronitromethane (Chloropicrin) | - | PCW |

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| :---: | :---: | :---: | :---: |
| 195 | Phosphorus oxychloride | - | PCW |
| 196 | Phosphorus trichloride | - | PCW |
| 197 | Phosphorus pentachloride | - | PCW |
| 198 | Trimethyl phosphite | - | PCW |
| 199 | Triethyl phosphite | - | PCW |
| 200 | Dimethyl phosphite | - | PCW |
| 201 | Diethyl phosphite | - | PCW |
| 202 | Sulfur monochloride | - | PCW |
| 203 | Sulfur dichloride | - | PCW |
| 204 | Thionyl chloride | - | PCW |
| 205 | Ethyldiethanolamine | - | PCW |
| 206 | Methyldiethanolamine | - | PCW |
| 207 | Triethanolamine | - | PCW |
| 208 | 1,1'-Dimethyl-4,4'-bipyridinium | - | PDSC |
| 209 | 1,1'-Dimethyl-4,4'-bipyridynium dimethylsulfate | - | PDSC |
| 210 | 1,1'-Dimethyl-4,4'-dipyridiniumdichloride | - | PDSC |
| 211 | 1,1-Dimethylhydrazine | - | PDSC |
| 212 | 1,3-Dichloropropane-2-ol | - | PDSC |
| 213 | Benzeneethanamine,4-fluoro-a-methyl,hydrochloride | - | PDSC |
| 214 | 1,3-Dinitro-4-chlorobenzene | - | PDSC |
| 215 | 1-Dodecylguanidinium acetate | - | PDSC |
| 216 | 2-Chloropyridine | - | PDSC |
| 217 | 2-Hydroxyethyl acrylate | - | PDSC |
| 218 | 2-Hydroxypropyl acrylate | - | PDSC |
| 219 | 2,2-Dimethyl-1,3-benzodioxol-4-yl-N- methylcarbamate | - | PDSC |
| 220 | 2,2-dimethylpropionyl chloride | - | PDSC |
| 221 | 2,3,5,6-Tetrafluoro-4- methylbenzyl (Z)- (1RS,3RS)-3-(2-chloro- 3,3,3-trifluoro-1-propenyl- 2,2-dimethylcyclopropane carboxylate | - | PDSC |
| 222 | 2,3-Dicyano-1,4-dithia-anthraquinone | - | PDSC |
| 223 | 2,4-Dinitro-6-(1- methylpropyl)-phenol | - | PDSC |
| 224 | 2-Diphenylacetyl-1,3- indandione | - | PDSC |
| 225 | 2-Mercaptoethanol | - | PDSC |
| 226 | 3-amino-1-propen | - | PDSC |
| 227 | 3-Chloropropane-1,2-diol | - | PDSC |
| 228 | Isophorone Diisocyanate | - | PDSC |
| 229 | 7-Bromo-6-chloro-3-[3- [(2R,3S)-3-hydroxy-2- piperidyl]-2-oxopropyl]-4(3H)-quinazolinone, | - | PDSC |
| 230 | ```N-Ethyl-methyl-(2-chloro- 4-methylmercaptophenyl)- thiophosphoramide``` | - | PDSC |
| 231 | O-Ethyl-O-(2- isopropoxycarbonylphenyl)- Nisopropylthiophosphoramide | - | PDSC |
| 232 | O-Ethyl S,S-dipropyl phosphorodithioate | - | PDSC |
| 233 | S,S-bis(1-methylpropyl) O-ethyl phosphorodithioate | - | PDSC |
| 234 | S-Methyl-N-[(methylcarbamoyl)-oxy]thioacetimidate | - | PDSC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 235 | Benzyl chloride | - | PDSC |
| 236 | Sodium azide | - | PDSC |
| 237 | Avermectin | - | PDSC |
| 238 | Allyl alcohol | - | PDSC |
| 239 | Dinoseb alkanolamine salts | - | PDSC |
| 240 | Ethylparanitrophenylthiono benzenephosphonate | - | PDSC |
| 241 | Octachlorotetrahydro methanophthalan | - | PDSC |
| 242 | Octamethyl pyrophosphoramide | - | PDSC |
| 243 | Tetramethyl orthosilicate | - | PDSC |
| 244 | Curare | - | PDSC |
| 245 | Crotonaldehyde | - | PDSC |
| 246 | Chloroacetaldehyde | - | PDSC |
| 247 | Methyl chloroacetate | - | PDSC |
| 248 | Phenyl chlorocarbonate | - | PDSC |
| 249 | Cobalt(II) oxide | - | PDSC |
| 250 | Sodium cyanide | - | PDSC |
| 251 | Hydrogen cyanide | - | PDSC |
| 252 | Diacetoxypropene | - | PDSC |
| 253 | Diethyl-(1,3- dithiocyclopentylidene)- thiophosphoramide | - | PDSC |
| 254 | Diethyl-4- methylsulfinylphenyl- thiophosphate | - | PDSC |
| 255 | Diethyl-S-(2-chloro-1- phthalimidoethyl)- dithiophosphate | - | PDSC |
| 256 | Diethyl-S-(ethylthioethyl)- dithiophosphate | - | PDSC |
| 257 | Diethyl-p-dimethylamionosulfonylphenylthiophosphate | - | PDSC |
| 258 | Diethyl paranitrophenyl thiophosphate | - | PDSC |
| 259 | Benzal Chloride | - | PDSC |
| 260 | Dinitro-o-cresol | - | PDSC |
| 261 | Dinitrocresol | - | PDSC |
| 262 | 4,6-dinitro-o-cresol sodium salt | - | PDSC |
| 263 | Dinitrophenol | - | PDSC |
| 264 | Dibutan-1-yl(dichloro)stannane | - | PDSC |
| 265 | Diborane | - | PDSC |
| 266 | Dimethyl-(isopropylthioethyl)- dithiophosphate | - | PDSC |
| 267 | Dimethyl-(diethylamido-1-chlorocrotonyl)-phosphate | - | PDSC |
| 268 | Dimethylethylmercapto ethylthiophosphate | - | PDSC |
| 269 | Dimethylparanitrophenyl thiophosphate | - | PDSC |
| 270 | Strychnine | - | PDSC |
| 271 | Strychnine nitrate | - | PDSC |
| 272 | Selenium | - | PDSC |
| 273 | Selenious acid | - | PDSC |
| 274 | Selenic acid | - | PDSC |
| 275 | Selenium hexafluoride | - | PDSC |
| 276 | Sodium selenite | - | PDSC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 277 | Sodium selenite | - | PDSC |
| 278 | Barium selenite | - | PDSC |
| 279 | Selenium dioxide | - | PDSC |
| 280 | Cadmium selenide sulfide | - | PDSC |
| 281 | Hydrogen selenide | - | PDSC |
| 282 | Ferrous selenide | - | PDSC |
| 283 | Thiosemicarbazide | - | PDSC |
| 284 | Tetraethylpyrophosphate | - | PDSC |
| 285 | Tetramethylammonium hydroxide | - | PDSC |
| 286 | Benzotrichloride | - | PDSC |
| 287 | Tributylamine | - | PDSC |
| 288 | Narasin | - | PDSC |
| 289 | Nicotine | - | PDSC |
| 290 | Nicotine,sulfate | - | PDSC |
| 291 | Nickel carbonyl | - | PDSC |
| 292 | Bis(4-isocyanatocyclohexan-1-yl)methane | - | PDSC |
| 293 | Hydrazine | - | PDSC |
| 294 | Fluorosulfonic acid | - | PDSC |
| 295 | Butyl 2,3-dihydro-2,2-dimethylbenzofuran-7 -yl N,N'-dimethyl-N,N'thiodicarbamate | - | PDSC |
| 296 | Ethyl bromoacetate | - | PDSC |
| 297 | Hexakis(beta,beta-dimethylphenethyl)-distannoxane | - | PDSC |
| 298 | Hexachloro-epoxy- octahydro-endo,endo- dimethanonaphthalene | - | PDSC |
| 299 | Hexachloro-hexahydro- methano-benzo- dioxathiepine oxide | - | PDSC |
| 300 | Hexachlorocyclopentadiene | - | PDSC |
| 301 | Benzenethiol | - | PDSC |
| 302 | Phosgene | - | PDSC |
| 303 | Methanesulfonyl chloride | - | PDSC |
| 304 | Methyl-N',N'-dimethyl-N-[(methylcarbamoyl)oxy]-1- thiooxamimidate | - | PDSC |
| 305 | Methylcyclohexyl-4- chlorophenylthiophosphate | - | PDSC |
| 306 | Methylphosphonic acid dichloride | - | PDSC |
| 307 | Methyl mercaptan | - | PDSC |
| 308 | Methylenebis(1- thiosemicarbazide | - | PDSC |
| 309 | Monofluoroacetate | - | PDSC |
| 310 | Fluoroacetamide | - | PDSC |
| 311 | Sodium fluoroacetate | - | PDSC |
| 312 | Isopropyl nitrite | - | PDSC |
| 313 | Butyl nitrite | - | PDSC |
| 314 | Benzensulfonyl chloride | - | PDSC |
| 315 | Phosphorus oxychloride | - | PDSC |
| 316 | Yellow phosphorus | - | PDSC |
| 317 | Phosphorus pentachloride | - | PDSC |
| 318 | Phosphorus trichloride | - | PDSC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 319 | Boron trichloride | - | PDSC |
| 320 | Phosphorous trifluoride | - | PDSC |
| 321 | Boron trifluoride | - | PDSC |
| 322 | Tetraalkyl lead | - | PDSC |
| 323 | Sulfur tetrafluoride | - | PDSC |
| 324 | Mercury | - | PDSC |
| 325 | Mercurous nitrate | - | PDSC |
| 326 | Phenylmercuric acetate | - | PDSC |
| 327 | Mercurous acetate | - | PDSC |
| 328 | Mercuric acetate | - | PDSC |
| 329 | Mercuric oxide | - | PDSC |
| 330 | Mercury(I) oxide | - | PDSC |
| 331 | Mercuric bromide | - | PDSC |
| 332 | Mercuric nitrate | - | PDSC |
| 333 | Mercuric iodide | - | PDSC |
| 334 | Mercuric chloride | - | PDSC |
| 335 | Mercury oxycyanide | - | PDSC |
| 336 | Mercuric cyanide | - | PDSC |
| 337 | Mercuric thiocyanate | - | PDSC |
| 338 | Thimerosal | - | PDSC |
| 339 | Sulfuryl fluoride | - | PDSC |
| 340 | Hydrogen fluoride | - | PDSC |
| 341 | Hydrofluoric acid | - | PDSC |
| 342 | Cadmium cyanide | - | PDSC |
| 343 | Potassium cyanide | - | PDSC |
| 344 | Calcium cyanide | - | PDSC |
| 345 | Potassium cobalt cyanide | - | PDSC |
| 346 | Potassium nickel cyanide | - | PDSC |
| 347 | Zinc cyanide | - | PDSC |
| 348 | Platinum-barium cyanide | - | PDSC |
| 349 | Silver cyanide | - | PDSC |
| 350 | Gold-potassium cyanide | - | PDSC |
| 351 | Mercuric cyanide | - | PDSC |
| 352 | Cuprous cyanide | - | PDSC |
| 353 | Potassium cuprocyanide | - | PDSC |
| 354 | Sodium cuprocyanide | - | PDSC |
| 355 | Lead cyanide | - | PDSC |
| 356 | Phosphorus trisulfide | - | PDSC |
| 357 | Phosphorus pentasulfide | - | PDSC |
| 358 | Hydrogen phosphide | - | PDSC |
| 359 | Tungsten hexafluoride | - | PDSC |
| 360 | Arsenic | - | PDSC |
| 361 | Tetraarsenic tetrasulfide | - | PDSC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 362 | Arsenic disulfide | - | PDSC |
| 363 | Arsenic trisulfide | - | PDSC |
| 364 | Lead arsenite | - | PDSC |
| 365 | Hydrogen arsenide | - | PDSC |
| 366 | Arsenic acid | - | PDSC |
| 367 | Potassium arsenate | - | PDSC |
| 368 | Calcium arsenate | - | PDSC |
| 369 | Sodium arsenate | - | PDSC |
| 370 | Manganese arsenate | - | PDSC |
| 371 | Zinc arsenate | - | PDSC |
| 372 | Lead arsenate | - | PDSC |
| 373 | Disodium hydrogenarsenate | - | PDSC |
| 374 | Calcium arsenate | - | PDSC |
| 375 | Ferric arsenate | - | PDSC |
| 376 | Copper arsenate | - | PDSC |
| 377 | Cacium arsenate fluoride | - | PDSC |
| 378 | Arsenic trifluoride | - | PDSC |
| 379 | Arsenic pentafluoride | - | PDSC |
| 380 | Lithium hexafluoroarsenate | - | PDSC |
| 381 | Arsenious acid | - | PDSC |
| 382 | Potassium arsenite | - | PDSC |
| 383 | Calcium arsenite | - | PDSC |
| 384 | Sodium arsenite | - | PDSC |
| 385 | Arsenic pentachloride | - | PDSC |
| 386 | Paris green | - | PDSC |
| 387 | Arsenic trichloride | - | PDSC |
| 388 | Arsenic pentaoxide | - | PDSC |
| 389 | Octamethyl pyrophosphoramide | - | PDSC |
| 390 | Diethyl paranitrophenyl thiophosphate | - | PDSC |
| 391 | Dimethyl-(diethylamido-1- chlorocrotonyl)-phosphate | - | PDSC |
| 392 | Dimethylethylmercapto ethylthiophosphate | - | PDSC |
| 393 | Dimethylparanitrophenyl thiophosphate | - | PDSC |
| 394 | Tetraethylpyrophosphate | - | PDSC |
| 395 | Monofluoroacetate | - | PDSC |
| 396 | Fluoroacetamide | - | PDSC |
| 397 | Sodium fluoroacetate | - | PDSC |
| 398 | Tetramethyl lead | - | PDSC |
| 399 | Tetraethyl lead | - | PDSC |
| 400 | Tetraalkyl lead | - | PDSC |
| 401 | Aluminium phosphide | - | PDSC |
| 402 | Polychlorinated dibenzofuran | - | POL |
| 403 | Polychlorinated dibenzo-para-dioxin | - | POL |
| 404 | Coplanar polychlorinated biphenyl | - | POL |

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| :---: | :---: | :---: | :---: |
| 405 | 3-acetoxy-6-dimethylamino-4,4-diphenylheptane (also called acethylmetadol) and its salts | - | NPC |
| 406 | a-3-acetoxy-6-dimethylamino-4,4-diphenylheptane (also called aacethylmetadol) and its salts | - | NPC |
| 407 | $\beta$-3-acetoxy-6-dimethylamino-4,4-diphenylheptane (also called $\beta$ acethylmetadol) and its salts | - | NPC |
| 408 | a-3-acetoxy-6-methylamino-4,4-diphenylheptane (also called noracymethadole) and its salts | - | NPC |
| 409 | 1-[2-(4-aminophenyl)ethyl]-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called anileridine) and its salts | - | NPC |
| 410 | N -allylnormorphine (also called nalorphine), its ester and their salts | - | NPC |
| 411 | 3-allyl-1-methyl-4-phenyl-4-(propionyloxy) piperidine (also called allylprodine) and its salts | - | NPC |
| 412 | ecgonine and its salts | - | NPC |
| 413 | 3-(N-ethyl-N-methylamino)-1,1-di-(2-thienyl)-1-buten (also called ethylmethylthiambutene) and its salts | - | NPC |
| 414 | a-3-ethyl-1-methyl-4-phenyl-4-(propionyloxy) piperidine (also called alphameprodine) and its salts | - | NPC |
| 415 | $\beta$-3-ethyl-1-methyl-4-phenyl-4-(propionyloxy) piperidine (also called betameprodine) and its salts | - | NPC |
| 416 | 2-(4-chlorobenzyl)-1-(diethylamino) ethyl-5-nitrobenzimidazole (also called clonitazene) and its salts | - | NPC |
| 417 | cocaine and other ecgonine ester and its salts | - | NPC |
| 418 | coca leaf | - | NPC |
| 419 | codeine, ethylmorphine, other morphine ether, and its salts | - | NPC |
| 420 | diacetylmorphine (also known as heroin), other morphine ester, and its salts | - | NPC |
| 421 | 1-(3-cyano-3,3-diphenylpropyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called diphenoxylate) and its salts | - | NPC |
| 422 | 4-cyano-2-dimethylamino-4,4-diphenylbutane (also called methadone intermediate) and its salts | - | NPC |
| 423 | 4-cyano-1-methyl-4-phenylpiperidine (also called pethidine intermediate $A$ ) and its salts | - | NPC |
| 424 | 1-(diethylamino)ethyl-2-(4-etoxybenzyl)-5-nitrobenzimidazole (also called etonitazene) and its salts | - | NPC |
| 425 | 3-diethylamino-1,1-di-(2-thienyl)-1-butene (also called diethylthienebutene) and its salts | - | NPC |
| 426 | dihydrocodeinone (also called hydrocodone), its ester, and their salts | - | NPC |
| 427 | dihydrocodeine, its ester, and their salts | - | NPC |
| 428 | dihydrodeoxy-morphine (also called desomorphine), its ester, and their salts | - | NPC |
| 429 | dihydrohydroxy-codeinone (also called oxycodone), its ester, and their salts | - | NPC |
| 430 | dihydrohydroxy-morphinone (also called oxyxmorphine) and its salts | - | NPC |
| 431 | dihydromorphine, its ester, and their salts | - | NPC |
| 432 | dihydromorphinone (also called hydromorphone), its ester, and their salts | - | NPC |
| 433 | 4,4-diphenyl-6-piperidinyl-3-heptanone (also called dipipanone) and its salts | - | NPC |
| 434 | (2-dimethylamino)ethyl-1-ethoxy-1,1-diphenylacetate (also called dimenoxadol) and its salts | - | NPC |
| 435 | 3-dimethylamino-1,1-di-(2-thienyl)-1-butene (also called dimethylthiambutene) and its salts | - | NPC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 436 | 6-dimethylamino-4,4-diphenyl-3-hexanone (also called normesadone) and its salts | - | NPC |
| 437 | 6-dimethylamino-4,4-diphenyl-3-heptanol (also called dimepheptanol) and its salts | - | NPC |
| 438 | a-6-dimethylamino-4,4-diphenyl-3-heptanol (also called a-methadol) and its salts | - | NPC |
| 439 | $\beta$-6-dimethylamino-4,4-diphenyl-3-heptanol (also called $\beta$-methadol) and its salts | - | NPC |
| 440 | 6-dimethylamino-4,4-diphenyl-3-heptanone (also called methodone) and its salts | - | NPC |
| 441 | 4-dimethylamino-3-methyl-1,2-diphenyl-2-(propionyloxy) butane (also called propoxyphine) and its salts | - | NPC |
| 442 | 6-dimethylamino-5-methyl-4,4-diphenyl-3-hexanone (also called isomethadone) and its salts | - | NPC |
| 443 | 1,3-dimethyl-4-phenyl-4-(propionyloxy)azacycloheputane (also called proheptazine) and its salts | - | NPC |
| 444 | a-1,3-dimethyl-4-phenyl-4-(propionyloxy) piperidine (also called aprodine) and its salts | - | NPC |
| 445 | $\beta$-1,3-dimethyl-4-phenyl-4-(propionyloxy) piperidine (also called $\beta$ prodine) and its salts | - | NPC |
| 446 | thebaine and its salts | - | NPC |
| 447 | 1,2,5-trimethyl-4-phenyl-4-(propionyloxy) piperidine (also called trimeperidine) and its salts | - | NPC |
| 448 | 6-nicotinylcodein (also called nicocodine) and its salts | - | NPC |
| 449 | normorphine (also called demethylmorphine), its ether, and their salts | - | NPC |
| 450 | 1-[2-(2-hydroxyetoxy)ethyl]-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called etoxyeridine) and its salts | - | NPC |
| 451 | 14-hydroxydihydromorphine (also called hydromorphinol) and its salts | - | NPC |
| 452 | 3-hydroxy-N-phenacylmorphinan (other than dextrorotatory one) and its salts | - | NPC |
| 453 | 1-(3-hydroxy-3-phenylpropyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called phenoperidine) and its salts | - | NPC |
| 454 | 4-(3-hydroxyphenyl)-1-methyl-4-piperidineethylketone (also called ketobemidone) and its salts | - | NPC |
| 455 | 4-(3-hydroxyphenyl)-1-methylpiperidine-4-carboxylic acid ethyl ester (also called hydroxypethidine) and its salts | - | NPC |
| 456 | 3-hydroxy-N-phenetylmorphinan (also called phenomorphan) and its salts | - | NPC |
| 457 | 3-hydroxy-N-methylmorphinan (other than dextrorotatory one) and its salts | - | NPC |
| 458 | 3-hydroxymorphinan (other than dextrorotatory one) and its salts | - | NPC |
| 459 | 4-phenyl-1-[2-(tetrahydrofurfuryloxy)ethyl]piperidine-4-carboxylic acid ethyl ester (also called furethidine) and its salts | - | NPC |
| 460 | 4-phenylpiperidine-4-carboxylic acid ethyl ester (also called pethidine intermediate $B$ ) and its salts | - | NPC |
| 461 | 4-phenyl-1-(3-pheminoaminopropyl)piperidine-4-carboxylic acid ethyl ester (also called piminodine) and its salts | - | NPC |
| 462 | 1,2,3,4,5,6-hexahydro-8-hydroxy-6,11-dimethyl-3-phenethyl-2,6-methano-3-benzazocine (also called phenazocine) and its salts | - | NPC |
| 463 | 1,2,3,4,5,6-hexahydro-8-hydroxy-3,6,11-trimethyl-2,6-methano-3penzazocine (also called metazocine) and its salts | - | NPC |
| 464 | 1-[2-(benzyloxy)ethyl]-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called benzethidine) and its salts | - | NPC |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 465 | 6-methyldihydromorphine (also called methyldihydromorphine) and its salts | - | NPC |
| 466 | methyldihydromorphinone (also called metopon), its ester, and their salts | - | NPC |
| 467 | 6-methyl- $\Delta$-6-dioxymorhine (also called metyldesorphine) and its salts | - | NPC |
| 468 | N -(1-methyl-2-piperidinoethyl)propyonanilide (also called phenanpromid) and its salts | - | NPC |
| 469 | 1-methyl-4-phenylpiperidine-4-carboxylic acid ester and its salts | - | NPC |
| 470 | N -[2-(methylphenetylamino)propyl]propyonanilide (also called diampromide) and its salts | - | NPC |
| 471 | [(3-methyl-4-morpholino-2,2-diphenyl)putyryl]piroridine and its salts | - | NPC |
| 472 | 3-methyl-4-morpholino-2,2-diphenyl butyrate (also called moramido intermediate) and its salts | - | NPC |
| 473 | 3-methoxy-N-methylmorphinan (other than dextrorotatory one) and its salts | - | NPC |
| 474 | morphine and its salts | - | NPC |
| 475 | morphine-N-oxydo, other pentavalent nitrogen morphine, and its derivatives | - | NPC |
| 476 | 1-(2-morpholinoethyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester (also called morphlidine) and its salts | - | NPC |
| 477 | 6-morpholino-4,4-diphenyl-3-heptanone (also called phenadoxone) and its salts | - | NPC |
| 478 | 4-morpholino-2,2-dipyenyl buturate ethyl ester (also called dioxaphetyl putyrate) and its salts | - | NPC |
| 479 | 3-O-acetyl-7,8-dihydro-7a-[1(R)-hydroxy-1-methylbutyl]-6-O-methyl-6,14-endo-ethenomorphine (also called acetorphine) and its salts | - | NPC |
| 480 | N -(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (also called APINACA N- (5-fluoropentyl) derivative) and its salts | - | NPC |
| 481 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamid (also called ADB-CHMINACA) and its salts | - | NPC |
| 482 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl )-1-butyl-1H-indazole-3carboxamide |  |  |
| 483 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (also called ADB-FUBINACA) and its salts | - | NPC |
| 484 | 3-(2-aminobutyl)indole (also called etryptamine) and its salts | - | NPC |
| 485 | 2-aminopropiophenone (also called cathinon) and its salts | - | NPC |
| 486 | 3-(2-aminopropyl)indole (also called AMT) and its salts | - | NPC |
| 487 | N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (also called AB-CHMINACA) and its salts | - | NPC |
| 488 | N - (1-amino-3-methyl-1-oxobutane-2-yl) -1- (4-fluorobenzyl) -1H-indazole-3-carboxamide and its salts |  | NPC |
| 489 | N -(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3carboxamide (also called AB-PINACA) and its salts | - | NPC |
| 490 | 2-ethylamino-1-phenylpropan-1-one (also called etocathinon) and its salts | - | NPC |
| 491 | 2- (Ethylamino)-1-phenylhexane-1-one and its salts |  | NPC |
| 492 | $\begin{aligned} & \text { 2- (ethylamino)-1-(4-methylphenyl)propan-1-one (also called 4-MEC) } \\ & \text { and its salts } \end{aligned}$ | - | NPC |
| 493 | 2-ethylamino-1- (3,4-methylenedioxyphenyl) propan-1-one (also called b $k$ - MDEA) and its salts | - | NPC |
| 494 | 2-ethylamino-1- (3,4-methylenedioxyphenyl) pentan-1-one (also called N -Ethylnorpentylone) and its salts | - | NPC |
| 495 | 2-Ethylamino-1-(3,4-methylenedioxyphenyl)pentan-1-one |  |  |

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| :---: | :---: | :---: | :---: |
| 496 | 2-(ethylamino)-2-(3-methoxyphenyl)cyclohexanone (also called Methoxetamine) and its salts | - | NPC |
| 497 | N -[1-[2-(4-ethyl-5-oxo-2-tetrazolin-1-yl)ethyl]-4-(methoxymethyl)-4piperidyl]propionanilide (also called alfentanil) and its salts | - | NPC |
| 498 | 4-ethyl-2,5-dimethoxy-a-methylphenethylamine (also called DOET) and its salts | - | NPC |
| 499 | 2-(4-ethylsulfanyl-2,5-dimethoxyphenyl)ethanamine (also called 2C-T- <br> 2) and its salts | - | NPC |
| 500 | N -ethyl-1-phenylcyclohexylamine (also called ethycyclidine) and its salts | - | NPC |
| 501 | N-ethyl-a-methyl-3,4-(methylenedioxy)phenethylamine (also called Nethyl MDA) and its salts | - | NPC |
| 502 | 2-(4-ethoxybenzyl)-5-nitro-1-[2-(pyrrolidin-1-yl)ethyl]benzimidazole |  |  |
| 503 | (5R)-4,5-epoxy-6-methoxy-17-methyl-6,7,8,14-tetradehydromorphinan-$3-\mathrm{ol}$ (also called oripavine) and its salts | - | NPC |
| 504 | Quinolin-8-yl=1-(5-fluoropentyl)-1H-indole-3-carboxylate (also called 5F-QUPIC) and its salts | - | NPC |
| 505 | 1- (4-Chloro-2,5-dimethoxyphenyl) propan-2-amine and its salts |  | NPC |
| 506 | 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (also called $2 \mathrm{C}-\mathrm{C}-\mathrm{NBOMe}$ ) and its salts | - | NPC |
| 507 | 1-(3-chlorophenyl)piperazine (also called 3CPP) and its salts | - | NPC |
| 508 | 2-(2-chlorophenyl)-2-(methylamino)cyclohexanone (also called ketamine) and its salts | - | NPC |
| 509 | 1- (4-chlorophenyl)-2-(methylamino) propan-1-one and its salts |  | NPC |
| 510 | 1-(3-cyano-3,3-diphenylpropyl)-4-(2-oxo-3-propionyl-1benzimidazolinyl)piperidine (also called bezitramide) and its salts | - | NPC |
| 511 | 1-(3-cyano-3,3-diphenylpropyl)-4-(1-piperidino)piperidine-4carboxamide (also called piritramide) and its salts | - | NPC |
| 512 | 1-(3-cyano-3,3-diphenylpropyl)-4-phenylpiperidine-4-carboxylic acid (also called difenoxin) and its salts | - | NPC |
| 513 | 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (also called CUMYL-4CN-BINACA) and its salts | - | NPC |
| 514 | $\mathrm{N}, \mathrm{N}$-diallyl-5-methoxytryptamine (also called 5-MeO-DALT) and its salts | - | NPC |
| 515 | 3-[2-(diisopropylamino)ethyl]-5-methoxyindole (also called 5-MeODIPT) and its salts | - | NPC |
| 516 | 1-(Diethylamino)ethyl-2-(4-isopropoxybenzyl)-5-nitrobenzimidazole | - | NPC |
| 517 | 3-[2-(diethylamino)ethyl]indole (also called DET) and its salts | - | NPC |
| 518 | 1-(2-diethylamino)ethyl-2-(4-ethoxybenzyl)benzimidazole |  |  |
| 519 | 1-(2-diethylamino)ethyl-5-nitro-2-(4-propoxybenzyl)benzimidazole |  |  |
| 520 | 1-(Diethylamino)ethyl-2-(4-methoxybenzyl)-5-nitrobenzimidazole |  |  |
| 521 | 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine (also called MT-45) and its salts | - | NPC |
| 522 | 3,4-dichloro-N- \{ [1-(dimethylamino)cyclohexyl] methyl\} benzamide (also called AH-7921) and its salts | - | NPC |
| 523 | 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also called U-47700) and its salts | - | NPC |
| 524 | cis-2-amino-4-methyl-5-phenyl-2-oxazoline (also called 4methylaminorex) and its salts | - | NPC |
| 525 | dihydrocodeinon-6-(carboxymethyl)oxime (also called codoxime) and its salts | - | NPC |
| 526 | 7-[(10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-yl)amino]heptanoic acid (also called amineptine) and its salts | - | NPC |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 527 | 7,8-dihydro-7-a-[1-(R)-hydroxy-1-methylbutyl]-6,14-endoethanotetrahydrooripavine(also called dihydroetorphine) and its salts | - | NPC |
| 528 | 7,8-dihydro-7a-[1(R)-hydroxy-1-methylbutyl]-6-O-methyl-6,14-endoethenomorphine (also called etorphine) and its salts | - | NPC |
| 529 | 1-(1,2-diphenylethyl)piperidine | - | NPC |
| 530 | 4,4-diphenyl-6-piperidino-3-hexanone (also called norpipanone) and its salts | - | NPC |
| 531 | 3-[2-(dimethylamino)ethyl]indole (also called DMT) and its salts | - | NPC |
| 532 | 3-[(2-dimethylamino)ethyl]-indol-4-yl phosphate ester (also called psilocybin) and its salts | - | NPC |
| 533 | 3-[2-(dimethylamino)ethyl]-indol-4-ol (also called psilocin) and its salts | - | NPC |
| 534 | 3-[(1R,2R)-3-(dimethylamino)-1-ethyl-2-methylpropyl]phenol (also called Tapentadol) and its salts | - | NPC |
| 535 | 3-(1,2-dimethylheptyl)-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H- dibenzo[b,d]pyran-1-ol (also called DMHP) and its salts | - | NPC |
| 536 | N,a-dimethyl-3,4-(methylenedioxy)phenethylamine (also called MDMA) and its salts | - | NPC |
| 537 | 2-(2,5-dimethoxy-4-isopropylsulfanylphenyl)ethaneamine (also called 2C-T-4) and its salts | - | NPC |
| 538 | 2,5-dimethoxy-4,a-dimethylphenethylamine (also called DOM) and its salts | - | NPC |
| 539 | 2,5-dimethoxy-4-(propylthio)phenethylamine and (also called 2C-T-7) and its salts | - | NPC |
| 540 | 2,5-dimethoxy-a-methylphenethylamine (also called DMA) and its salts | - | NPC |
| 541 | 3,4-dimethoxy-17-methylmorphinan-6 $\beta, 14$-diol (also called drotebanol) and its salts | - | NPC |
| 542 | N-[1-[2-(2-thienyl)ethyl]-4-piperidyl]propionanilide (also called thiofentanyl) and its salts | - | NPC |
| 543 | 1-[1-(2-thienyl)cyclohexyl]piperidine (also called thenocyclidine) and its salts | - | NPC |
| 544 | 6a,7,8,9-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol (also called $\Delta 10$-tetrahydrocannabinol) and its salts | - | NPC |
| 545 | 6a,7,8,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1ol (also called $\Delta 9$-tetrahydrocannabinol)(limited to those obtainable from causing a chemical reaction other than decomposition reaction (excepting those necessary to purify 6a,7,8,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol contained in hemp stipulated in Article 1 of the Marijuana Control Act (Act No. 124 of 1948) (hereinafter simply referred to as "hemp") and its products)) and its salts | - | NPC |
| 546 | 6a,7,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol (also called $\Delta 8$-tetrahydrocannabinol) (limited to those obtainable from causing a chemical reaction other than decomposition reaction (excepting those necessary to purify 6a,7,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol contained in hemp and its products)) and its salts | - | NPC |
| 547 | 6a,9,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol (also called $\Delta 7$-tetrahydrocannabinol) and its salts | - | NPC |
| 548 | 7,8,9,10-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol (also called $\Delta 6 \mathrm{a}(10 \mathrm{a})$-tetrahydrocannabinol) and its salts | - | NPC |
| 549 | 8,9,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1ol (also called $\Delta 6 \mathrm{a}(7)$-tetrahydrocannabinol) and its salts | - | NPC |
| 550 | trans-2-dimethylamino-1-phenyl-3-cyclohexene-1-carboxylic acid ethyl ester (also called tilidine) and its salts | - | NPC |

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| :---: | :---: | :---: | :---: |
| 551 | 1-(3-trifluoromethylphenyl)piperazine (also called T F M P P ) and its salts | - | NPC |
| 552 | 3,4,5-trimethoxyphenethylamine (also called mescaline) and its salts | - | NPC |
| 553 | 2,4,5-trimethoxy-a-methylphenethylamine (also called TMA-2) and its salts | - | NPC |
| 554 | 3,4,5-trimethoxy-a-methylphenethylamine (also called TMA) and its salts | - | NPC |
| 555 | 1-naphthalenyl(1-pentyl-1H-indol-3-yl)methanone (also called JWH018) and its salts | - | NPC |
| 556 | N -[1-( $\beta$-hydroxyphenethyl)-4-piperidyl]propionanilide (also called $\beta$ hydroxyfentanyl) and its salts | - | NPC |
| 557 | N-[1-( $\beta$-hydroxyphenethyl)-3-methyl-4-piperidyl]propionanilide (also called $\beta$-hydroxy-3-methylfentanyl) and its salts | - | NPC |
| 558 | (1RS, 3SR) -3- [2-hydroxy-4- (2-methylnonan-2-yl) phenyl] cyclohexane-1-ol (also called Cannabicyclohexal) and its salts | - | NPC |
| 559 | 4-hydroxybutyric acid (also called GHB) and its salts | - | NPC |
| 560 | 1-(1-phenylcyclohexyl)piperidine (also called phencyclidine) and its salts | - | NPC |
| 561 | 1-(1-phenylcyclohexyl)pyrrolidine (also called rolicyclidine) and its salts | - | NPC |
| 562 | 2-phenyl-2- (piperidin-2-yl) acetic acid ethyl ester (also called ethylphenidate) and its salts | - | NPC |
| 563 | 1-Phenyl-2- (pyrrolidin-1-yl) hexane-1-one and its salts |  | NPC |
| 564 | 1-phenyl-2- (pyrrolidin-1-yl) pentan-1-one (also called a-P V P) and its salts | - | NPC |
| 565 | N-(1-phenethyl-4-piperidyl)propionanilide (also called fentanyl) and its salts | - | NPC |
| 566 | N -(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (also called acrylfentanyl) and its salts | - | NPC |
| 567 | N -(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also called acetylfentanyl) and its salts | - | NPC |
| 568 | N -(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (also called cyclopropylfentanyl) and its salts | - | NPC |
| 569 | N -(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (also called THF-F) and its salts | - | NPC |
| 570 | (E) -N - (1-phenethylpiperidin-4-yl) - N -phenylbuta-2-enamid and salts thereof |  | NPC |
| 571 | N -(1-phenethylpiperidin-4-yl)-N-phenylbutanamide (also called butyrlfentanyl) and its salts | - | NPC |
| 572 | N -(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (also called furanylfentanyl) and its salts | - | NPC |
| 573 | N - (1-phenethylpiperidin-4-yl) - N -phenylpentaneamide and its salts |  | NPC |
| 574 | 1-phenethyl-4-phenyl-4-piperidinol acetate (also called PEPAP) and its salts | - | NPC |
| 575 | (1-butyl-1H-indol-3-yl)(naphthalen-1-yl)methanon (also called J WH073 ) and its salts | - | NPC |
| 576 | N -(4-fluorophenyl)- N -(1-phenethylpiperidin-4-yl)butanamide (also called 4-FBF) and its salts | - | NPC |
| 577 | N -(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propanamide (also called 2-fluorofentanyl) and its salts | - | NPC |
| 578 | 1-(4-fluorophenyl)propan-2-amine (also called 4-FA) and its salts | - | NPC |
| 579 | N-(4-fluorophenyl)-2-methyl-N-(1-phenethylpiperidin-4yl)propanamide (also called 4-FIBF) and its salts | - | NPC |
| 580 | N -(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl) acetamide (also called Ocfentanil) and its salts | - | NPC |

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| :---: | :---: | :---: | :---: |
| 581 | 4-fluoro-N-(1-phenethyl-4-piperidyl)propionanilide (also called $\rho$ fluorofentanyl) and its salts | - | NPC |
| 582 | [1-(5-fluoropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropan-1yl)methanone (also called XLR-11) and its salts | - | NPC |
| 583 | [1-(5-fluoropentyl)-1H-indol-3-yl](naphthalen-1-yl)methanone (also called AM2201) and its salts | - | NPC |
| 584 | [1-(5-fluoropentyl)-1H-indol-3-yl](4-methylnaphthalen-1-yl)methanone (also called MAM-2201) and its salts | - | NPC |
| 585 | 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (also called 25B-NBOMe) and its salts | - | NPC |
| 586 | 4-bromo-2,5-dimethoxyphenethylamine (also called 2C-B) and its salts | - | NPC |
| 587 | 4-bromo-2,5-dimethoxy-a-methylphenethylamine (also called brolamfetamine) and its salts | - | NPC |
| 588 | 1-\{1-\{1-[1-(4-bromophenyl)ethyl]piperidin-4-yl\}-1,3-dihydro-2H-benzo[d]imidazol-2-one |  |  |
| 589 | 6a,7,8,9,10,10a-hexahydro-6,6-dimethyl-9-methylene-3-pentyl-6H-dibenzo[b,d]pyran-1-ol(also called $\Delta 9$ (11)-tetrahydrocannabinol) and its salts | - | NPC |
| 590 | 3-hexyl-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran-1-ol (also called parahexyl) and its salts | - | NPC |
| 591 | 1-benzylpiperadine (also called B Z P) and its salts | - | NPC |
| 592 | (1-pentyl-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropan-1yl)methanone (also called UR-144) and its salts | - | NPC |
| 593 | 5-Pentyl-2-(2-phenylpropan-2-yl)-2,5-dihydro-1H-pyrido[4,3-b]indol-1one | - | NPC |
| 594 | 2-(methylamino)-1-phenylpropan-1-one (also called methcathinone) and its salts | - | NPC |
| 595 | 2-(methylamino)-1-phenylpentan-1-one (also called pentedrone) and its salts | - | NPC |
| 596 | 2-(methylamino)-1-(3-methylphenyl)propan-1-one |  |  |
| 597 | 2-(methylamino)-1-(4-methylphenyl)propan-1-one (also called 4methylmethcathinone ) and its salts | - | NPC |
| 598 | 2-methylamino-1-(3,4-methylenedioxyphenyl)propan-1-one (also called metilon) and its salts | - | NPC |
| 599 | N -methyl-a-ethyl-3,4-(methylenedioxy)phenethylamine (also called MBDB) and its salts | - | NPC |
| 600 | methyl=2-[1-(cyclohexylmethyl)-1H-indole-3-carboxamide]-3,3dimethylbutanoate (also called MDMB-CHIMICA) and its salts | - | NPC |
| 601 | Methyl=3,3-dimethyl-2-[1-(penta-4-en-1-yl)-1H-indazole-3carboxamido]butanoate | - | NPC |
| 602 | N-[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]propionanilide (also called a-methylthiofentanyl) and its salts | - | NPC |
| 603 | N -[3-methyl-1-[2-(2-thienyl)ethyl]-4-piperidyl]propionanilide (also called 3-methylthiofentanyl) and its salts | - | NPC |
| 604 | N -methyl-1-(thiophen-2-yl)propan-2-amine (also called methiopropamine) and its salts | - | NPC |
| 605 | (4-methylnaphthalen-1-yl)(1-pentyl-1H-indol-3-yl)methanone (also called JWH-122) and its salts | - | NPC |
| 606 | N -[1-methyl-2-(piperidinoethyl)]-N-2-pyridylpropionamide (also called propiram) and its salts | - | NPC |
| 607 | 1-methyl-4-phenyl-4-piperidinol propionate (also called MPPP) and its salts | - | NPC |
| 608 | 1-methyl-4-phenylpiperidine-4-carboxylic acid (also called pethidine intermediate C) and its salts | - | NPC |
| 609 | 4-Methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one |  |  |

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| :---: | :---: | :---: | :---: |
| 610 | 1-[2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl]butan-1-one |  |  |
| 611 | N -[1-(a-methylphenethyl)-4-piperidyl]acetanilide (also called acetyl-amethylfentanyl) and its salts | - | NPC |
| 612 | N -[1-(a-methylphenethyl)-4-piperidyl]propionanilide (also called amethylfentanyl) and its salts | - | NPC |
| 613 | N-(3-methyl-1-phenethyl-4-piperidyl)propionanilide (also called 3methylfentanyl) and its salts | - | NPC |
| 614 | methyl=1-phenethyl-4-(N-phenylpropanamide)piperidine-4carboxylate (also called carfentanil) and its salts | - | NPC |
| 615 | Methyl = 2- [1- (4-fluorobutyl) -1H-indazole-3-carboxamide] -3,3dimethylbutanoate and its salts |  | NPC |
| 616 | methyl=2-[1-(4-fluorobenzyl)-1H-indazole-3-carboxamide]-3methylbutanoate (also called FUB-AMB) and its salts | - | NPC |
| 617 | methyl=2-[1- (5-fluoropentyl)-1H-indazole-3-carboxamide]-3,3dimethylbutanoate (also called 5F-ADB) and its salts | - | NPC |
| 618 | Methyl = 2- [1- (5-fluoropentyl) -1H-indazole-3-carboxamide] -3methylbutanoate and its salts |  | NPC |
| 619 | Methyl = 2- [1- (5-fluoropentyl) -1H-indole-3-carboxamide] -3,3dimethylbutanoate and its salts |  | NPC |
| 620 | a-methyl-4-methylthiophenethylamine (also called 4-MTA) and its salts | - | NPC |
| 621 | 4-methyl-5-(4-methylphenyl)-4,5-dihydrooxazol-2-amine (also called 4,4'-DMAR) and its salts | - | NPC |
| 622 | N-methyl-N-(1-(3,4-methylenedioxyphenyl)propan-2-yl)hydroxylamine (also called N-hydroxy MDMAN-OH MDMA) and its salts | - | NPC |
| 623 | a-methyl-3,4-(methylenedioxy)phenethylamine (also called MDA) and its salts | - | NPC |
| 624 | N-[a-methyl-3,4-(methylenedioxy)phenethyl]hydroxylamine (also called N -hydroxy MDA) and its salts | - | NPC |
| 625 | 1-(3,4-methylenedioxyphenyl) -2-(pyrrolidin-1-yl)pentan-1-one (also called MDPV) and its salts | - | NPC |
| 626 | 1-(2-methoxycarbonylethyl)-4-(phenylpropionylamino)piperidine-4carboxylic acid methyl ester (also called remifentanil) and its salts | - | NPC |
| 627 | 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine | - | NPC |
| 628 | 1-(4-methoxyphenyl)-N-methylpropan-2-amine (also called PMMA) and its salts | - | NPC |
| 629 | 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also called methoxyacetylfentanyl) and its salts | - | NPC |
| 630 | N -[4-(methoxymethyl)-1-[2-(2-thienyl)ethyl]-4-piperidyl]propionanilide (also called sufentanil) and its salts | - | NPC |
| 631 | 4-methoxy-a-methylphenethylamine (also called PMA) and its salts | - | NPC |
| 632 | 3-methoxy-a-methyl-4,5-(methylenedioxy)phenethylamine (also called MMDA) and its salts | - | NPC |
| 633 | 2-(4-iodo-2,5-dimethoxyphenyl)ethanamine (also called 2C-I) and its salts | - | NPC |
| 634 | 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (also called 251-NBOMe) and its salts | - | NPC |
| 635 | lysergic acid diethylamide (also called lysergide) and its salts | - | NPC |
| 636 | (Opium) Poppy | - | NPC |
| 637 | Papavel Somniferum El | - | NPC |
| 638 | Papaver Setighelm Dc | - | NPC |
| 639 | carbon dioxide | - | PGWC |
| 640 | methane | - | PGWC |
| 641 | Nitrous oxide | - | PGWC |

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| :---: | :---: | :---: | :---: |
| 642 | Trifluoromethane (aka HFC-23) | - | PGWC |
| 643 | Difluoromethane (aka HFC-32) | - | PGWC |
| 644 | Fluoromethane (aka HFC-41) | - | PGWC |
| 645 | 1.1.1.2.2-Pentafluoroethane (also known as HFC-125) | - | PGWC |
| 646 | 1.1.2.2-Tetrafluoroethane (also known as HFC-134) | - | PGWC |
| 647 | 1.1.1.2-Tetrafluoroethane (also known as HFC-134a) | - | PGWC |
| 648 | 1.1.2-Trifluoroethane (aka HFC-143) | - | PGWC |
| 649 | 1.1.1-Trifluoroethane (also known as HFC-143a) | - | PGWC |
| 650 | 1.2-Difluoroethane (aka HFC-152) | - | PGWC |
| 651 | 1.1-Difluoroethane (aka HFC-152a) | - | PGWC |
| 652 | Fluoroethane (aka HFC-161) | - | PGWC |
| 653 | 1.1.1.2.3.3.3-Heptafluoropropane (also known as HFC-227ea) | - | PGWC |
| 654 | 1.1.1.3.3.3-Hexafluoropropane (aka HFC-236fa) | - | PGWC |
| 655 | 1.1.1.2.3.3-Hexafluoropropane (also known as HFC-236ea) | - | PGWC |
| 656 | 1.1.1.2.2.3-Hexafluoropropane (also known as HFC-236 cb) | - | PGWC |
| 657 | 1.1.2.2.3-Pentafluoropropane (also known as HFC-245ca) | - | PGWC |
| 658 | 1.1.1.3.3-Pentafluoropropane (also known as HFC-245fa) | - | PGWC |
| 659 | 1.1.1.3.3-Pentafluorobutane (also known as HFC-365mfc) | - | PGWC |
| 660 | 1.1.1.2.3.4.4.5.5.5-Decafluoropentane (also known as HFC-43-10mee) | - | PGWC |
| 661 | Perfluoromethane (aka PFC-14) | - | PGWC |
| 662 | Perfluoroethane (aka PFC-116) | - | PGWC |
| 663 | Perfluoropropane (also known as PFC-218) | - | PGWC |
| 664 | Perfluorocyclopropane | - | PGWC |
| 665 | Perfluorobutane (also known as PFC-31-10) | - | PGWC |
| 666 | Perfluorocyclobutane (also known as PFC-c318) | - | PGWC |
| 667 | Perfluoropentane (also known as PFC-41-12) | - | PGWC |
| 668 | Perfluorohexane (also known as PFC-51-14) | - | PGWC |
| 669 | Perfluorodecalin (also known as PFC-91-18) | - | PGWC |
| 670 | Sulfur hexafluoride | - | PGWC |
| 671 | Nitrogen trifluoride | - | PGWC |
| 672 | Isobutyl nitrite | - | SQE |
| 673 | Isopropyl nitrite | - | SQE |
| 674 | Isopentyl nitrite | - | SQE |
| 675 | tert-Butyl nitrite | - | SQE |
| 676 | cyclohexyl nitrite | - | SQE |
| 677 | Butyl nitrite | - | SQE |
| 678 | 4-acetyl-N, N-diethyl-7-methyl-4,6,6a, 7,8,9-hexahydroindolo [4,3-fg] quinoline-9-carboxamide and salts thereof | - | SQE |
| 679 | 4-Acetoxy-N-Isopropyl-N-Methyltryptamine and its salts | - | SQE |
| 680 | 4-Acetoxy-N-ethyl-N-methyltryptamine and its salts | - | SQE |
| 681 | 4-Acetyloxy-N,N-diallyltryptamine and its salts | - | SQE |
| 682 | 4-Acetoxy-N,N-diisopropyltryptamine and its salts | - | SQE |
| 683 | 4-Acetoxy-N,N-diethyltryptamine and its salts | - | SQE |

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| :---: | :---: | :---: | :---: |
| 684 | 4-Acetoxy-N,N-dimethyltryptamine and its salts | - | SQE |
| 685 | N - (1-adamantyl) -1- (5-chloropentyl) -1H-indazole-3carboxamide and its salts | - | SQE |
| 686 | N -(1-Adamantyl)-1-(cyclohexylmethyl)-1H-indazole-3carboxamide | - | SQE |
| 687 | N -(1-adamanty)-1-[(tetrahydro-2H-pyran-4 (tetrahydro-2H-pyran-4-yl)methyl]-1H-indazole-3-carboxamide |  |  |
| 688 | N - (2-adamantyl) -1 -[(tetrahydro-2H-pyran-4-yl) methyl] -1H-indazole-3-carboxamide and its salts | - | SQE |
| 689 | N -(1-adamantyl)-1-(4-fluorobutyl)-1 H -indazol-3-carboxamide |  |  |
| 690 | FUB-APINACA |  | SQE |
| 691 | N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide and its salts | - | SQE |
| 692 | N -(1-adamantyl)-1-pentylindazole-3-carboxamide and its salts | - | SQE |
| 693 | 1 -adamantyl $=1$-pentyl-1H-indazole-3-carboxylate and its salts | - | SQE |
| 694 | 1-adamantyl-(1-pentylindol-3-yl)methanone and its salts | - | SQE |
| 695 | N -(1-adamantyl)-1-pentylindole-3-carboxamide and its salts | - | SQE |
| 696 | 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3yl]methanone and its salts | - | SQE |
| 697 | N -[(2S)-1-amino-1-oxo-3-phenylpropan-2-yl]-1-(cyclohexylmethyl)indazole-3-carboxamide and its salts and its salts | - | SQE |
| 698 | N -(1-amino-1-oxo-3-phenylpropan-2-yl)-1-butyl-1H-indazole-3carboxamide |  |  |
| 699 | $\mathrm{N}-[(2 \mathrm{~S})-1$-amino-1-oxo-3-phenylpropan-2-yl]-1-[(4-fluorophenyl)methyl]indazole-3-carboxamide and its salts | - | SQE |
| 700 | N - (1-Amino-1-oxo-3-phenylpropan-2-yl) -1- (5-fluoropentyl) 1 H -indazole-3-carboxamide and its salts | - | SQE |
| 701 | N - (1-Amino-1-oxo-3-phenylpropan-2-yl) -1- (5-fluoropentyl) 1 H -indole-3-carboxamide and its salts | - | SQE |
| 702 | N - (1-amino-3,3-dimethyl-1-oxobutan-2-yl) -1(cyclohexylmethyl) -1 H -indole-3-carboxamide and its salts | - | SQE |
| 703 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide and its salts $\mathrm{N}-[(2 S)-1$-amino-3,3-dimethyl-1-oxobutan-2-yl]-1-(5-fluoropentyl)indazole-3-carboxamide and its salts | - | SQE |
| 704 | N -(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentan-1- <br> y ) - 1 H -indole- 3 -carboxamide and its salts <br> N -((1S)-1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5- <br> fluoropentan- 1 -yl)-1 H -indole-3-carboxamide and its salts | - | SQE |
| 705 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-hexyl-1H-indazole-3-carboxamide |  |  |
| 706 | N -(1-amino-3,3-dimethyl-1-oxobutan-2-yl ) -1-benzyl-1H-indazole-3-carboxamide |  |  |
| 707 | N -(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide and its salts | - | SQE |
| 708 | N -(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole3 -carboxamide and its salts | - | SQE |
| 709 | 2-amino-1-(4-bromo-2,5-dimethoxyphenyl)ethanone and its salts | - | SQE |


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| :---: | :---: | :---: | :---: |
| 710 | N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-chloropentyl)indazole-3-carboxamide and its salts N -[(2S)-1-amino-3-methyl-1-oxobutan-2-yl]-1-(5-chloropentyl)indazole-3-carboxamide and its salts | - | SQE |
| 711 | $\mathrm{N}-[(2 \mathrm{~S})-1$-amino-3-methyl-1-oxobutan-2-yl]-1-[(2-fluorophenyl)methyl]indazole-3-carboxamide and its salts | - | SQE |
| 712 | N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide and its salts N -[(2S)-1-amino-3-methyl-1-oxobutan-2-yl]-1-(5-fluoropentyl)indazole-3-carboxamide and its salts | - | SQE |
| 713 | N -(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indole-3-carboxamide and its salts <br> N -[(2S)-1-amino-3-methyl-1-oxobutan-2-yl]-1-(5- <br> fluoropentyl)indole-3-carboxamide and its salts | - | SQE |
| 714 | 2-(3,5-dimethoxy-4-prop-2-enoxyphenyl)ethanamine 2-(3,5-dimethoxy-4-prop-2enoxyphenyl)ethanamine;hydrochloride | - | SQE |
| 715 | 7-allyl-N, N-diethyl-4,6,6a, 7,8,9-hexahydroindolo [4,3-fg] quinoline- 9 -carboxamide and salts thereof | - | SQE |
| 716 | 2-(Isopropylamino)-2-(3-methoxyphenyl)cyclohexanone |  |  |
| 717 | ```1-(2,5-dimethoxy-4-propan-2-ylsulfanylphenyl)propan-2-amine and its salts 1-(4-isopropylsulfanyl-2,5-dimethoxyphenyl)propan-2-amine and its salts``` | - | SQE |
| 718 | n-methyl-n-isopropyltryptamine and its salts | - | SQE |
| 719 | 5-Methoxy-N-methyl-N-isopropyltryptamine and its salts N-Isopropyl-N-methyl-5-methoxytryptamine hydrochloride | - | SQE |
| 720 | Nitrous oxide | - | SQE |
| 721 | Indan-2-amine and its salts 2-Aminoindan hydrochloride | - | SQE |
| 722 | 1- (Indan-5-yl) -2- (pyrrolidin-1-yl) butan-1-one and its salts | - | SQE |
| 723 | 1- (Indan-5-yl)-2-(pyrrolidin-1-yl) hexane-1-one and its salts | - | SQE |
| 724 | (1H-Indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone and its salts | - | SQE |
| 725 | 1-(1H-indol-5-yl)propan-2-amine and its salts | - | SQE |
| 726 | 2- (ethylamino) -1- (indan-5-yl) butan-1-one and its salts | - | SQE |
| 727 | 2- (ethylamino) -1- (indan-5-yl) pentan-1-one and its salts | - | SQE |
| 728 | 3- [1- (ethylamino) cyclohexyl] phenol and salts thereof | - | SQE |
| 729 | $\begin{aligned} & \text { 2-(Ethylamino)-2-(2-thienyl)cyclohexanone and its salts } \\ & \text { 2-(ethylamino)-2-thiophen-2-ylcyclohexan-1-one;hydrochloride } \end{aligned}$ | - | SQE |
| 730 | 2-(ethylamino) -2-phenylcyclohexanone and its salts 2-(ethylamino)-2-phenylcyclohexan-1-one;hydrochloride | - | SQE |
| 731 | 2-(Ethylamino)-2-(3-fluorophenyl)cyclohexanone |  |  |
| 732 | 2-(Ethylamino)-2-(3-methylphenyl)cyclohexanone |  |  |
| 733 | N -Ethyl-N-isoprpyl-5-methoxy-tryptamine and its salts | - | SQE |
| 734 | N -ethyl-1,2-diphenylethanamine and its salts N-ethyl-1,2-diphenylethanamine;hydrochloride | - | SQE |
| 735 | 2-(4-ethyl-2,5-dimethoxyphenyl)ethanamine and its salts 2,5-Dimethoxy-4-ethylphenethylamine hydrochloride | - | SQE |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 736 | 2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2methoxyphenyl)methyl]ethanamine and its salts 2-(4-Ethyl-2,5-dimethoxyphenyl)-N-[(2methoxyphenyl)methyl]ethanamine;hydrochloride | - | SQE |
| 737 | ```2-[[2-(4-Ethyl-2,5-dimethoxyphenyl)ethylamino]methyl]phenol and its salts 2-[[2-(4-Ethyl-2,5- dimethoxyphenyl)ethylamino]methyl]phenol;hydrochloride``` | - | SQE |
| 738 | ```1-(4-ethylsulfanyl-2,5-dimethoxyphenyl)propan-2-amine and its salts 1-(4-ethylsulfanyl-2,5-dimethoxyphenyl)propan-2- amine;hydrochloride``` | - | SQE |
| 739 | 4-Hydroxy-N-methyl-N-ethyltryptamine and its salts | - | SQE |
| 740 | 1- (4-Ethylphenyl) -N- (2-methoxybenzyl) propane-2-amine and its salts | - | SQE |
| 741 | Ethyl 2-(4-fluorophenyl)-2-(piperidin-2-yl)acetate | - | SQE |
| 742 | N -Ethyl-1-(2-fluorophenyl)propan-2-amine and its salts N-Ethyl-1-(2-fluorophenyl)propan-2-amine;hydrochloride | - | SQE |
| 743 | N-Ethyl-1-(3-fluorophenyl)propan-2-amine and its salts N-Ethyl-1-(3-fluorophenyl)propan-2-amine;hydrochloride | - | SQE |
| 744 | N-Ethyl-1-(4-fluorophenyl)propan-2-amine and its salts N-Ethyl-1-(4-fluorophenyl)propan-2-amine;hydrochloride | - | SQE |
| 745 | 3-ethyl-2-(3-fluorophenyl) morpholine and its salts | - | SQE |
| 746 | Ethyl 2- [1- (4-fluorobenzyl) -1H-indazole-3-carboxamide] -3methylbutanoate and its salts | - | SQE |
| 747 | Ethyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3dimethylbutanoate | - | SQE |
| 748 | Ethyl 2- [1- (5-fluoropentyl) -1H-indazole-3-carboxamide] -3methylbutanoate and its salts | - | SQE |
| 749 | Ethyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3dimethylbutanoate | - | SQE |
| 750 | Ethyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3methylbutanoate | - | SQE |
| 751 | 3-\{2-[Ethyl(propyl)amino]ethyl\}-1H-indol-4-yl acetate | - | SQE |
| 752 | N-ethyl-N-methyltryptamine |  |  |
| 753 | N -ethyl- N -[2-(5-methoxy-1H-indol-3-yl)ethyl]propan-1-amine and its salts | - | SQE |
| 754 | 3-MeO-PCE;3-Methoxyeticyclidine and its salts N-Ethyl-1-(3-methoxyphenyl)cyclohexan-1-amine and its salts N -Ethyl-1-(3-methoxyphenyl)cyclohexan-1-amine;hydrochloride | - | SQE |
| 755 | N-ethyl-1-(4-methoxyphenyl)propan-2-amine and its salts N-ethyl-1-(4-methoxyphenyl)propan-2-amine;hydrochloride | - | SQE |
| 756 | 2-(4-ethoxy-3,5-dimethoxyphenyl)ethanamine and its salts 2-(4-ethoxy-3,5-dimethoxyphenyl)ethanamine;hydrochloride | - | SQE |
| 757 | 1-(4-ethoxy-3,5-dimethoxyphenyl)propan-2-amine and its salts | - | SQE |
| 758 | 2-(4-ethoxybenzyl)-5-nitro-1-[2-(piperidin-1yl)ethyl]benzimidazole |  |  |
| 759 | 2-(1-oxo-1-phenylpropan-2-yl)isoindole-1,3-dione and its salts | - | SQE |
| 760 | quinolin-8-yl 1-(cyclohexylmethyl)indole-3-carboxylate and its salts | - | SQE |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 761 | Quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate | - | SQE |
| 762 | Quinolin-8-yl 1- (4-fluorobenzyl)-1H-indazole-3-carboxylate | - | SQE |
| 763 | quinolin-8-yl 1-[(4-fluorophenyl)methyl]indole-3-carboxylate and its salts | - | SQE |
| 764 | quinolin-8-yl 1-(5-fluoropentyl)indazole-3-carboxylate and its salts | - | SQE |
| 765 | quinolin-8-yl 1-pentylindazole-3-carboxylate and its salts | - | SQE |
| 766 | quinolin-8-yl 1-pentylindole-3-carboxylate and its salts | - | SQE |
| 767 | 5-chloro-3-ethyl-N-[2-(4-piperidin-1-ylphenyl)ethyl]-1H-indole-2-carboxamide and its salts | - | SQE |
| 768 | 2-(4-chloro-2,5-dimethoxyphenyl)ethanamine and its salts <br> 2-(4-chloro-2,5-dimethoxyphenyl)ethanamine;hydrochloride | - | SQE |
| 769 | 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(3,4,5trimethoxyphenyl)methyl]ethanamine and its salts | - | SQE |
| 770 | 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2fluorophenyl)methyl]ethanamine and its salts 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2fluorophenyl)methyl]ethanamine;hydrochloride | - | SQE |
| 771 | 2-[[2-(4-chloro-2,5- <br> dimethoxyphenyl)ethylamino]methyl]phenol and its salts 2-[[2-(4-chloro-2,5- <br> dimethoxyphenyl)ethylamino]methyl]phenol;hydrochloride |  | SQE |
| 772 | 1-(4-chloro-2,5-dimethoxyphenyl)propan-2-amine and its salts 1-(4-chloro-2,5-dimethoxyphenyl)propan-2amine;hydrochloride | - | SQE |
| 773 | 2- (2-Chlorophenyl) -1- (1-pentyl-1H-indole-3-yl) etanone and its salts | - | SQE |
| 774 | 1-(4-chlorophenyl)propan-2-amine and its salts (2R)-1-(4-chlorophenyl)propan-2-amine and its salts (2S)-1-(4-chlorophenyl)propan-2-amine;hydrochloride 1-(4-chlorophenyl)propan-2-amine;hydrochloride (2R)-1-(4-chlorophenyl)propan-2-amine;hydrochloride | - | SQE |
| 775 | 4-Chloroisobutyrylfentanyl and its salts <br> 1- (4-Chlorophenyl) - N -methylpropane-2-amine and its salts | - | SQE |
| 776 | 2-(3-chlorophenyl)-3-methylmorpholine |  |  |
| 777 | Salvinorin A | - | SQE |
| 778 | 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)pyrrolo[2,3-b]pyridine-3-carboxamide and its salts | - | SQE |
| 779 | $\mathrm{N}, \mathrm{N}$-Diisopropyltryptamine and its salts N -[2-(1H-indol-3-yl)ethyl]-N-propan-2-ylpropan-2-amine and its salts | - | SQE |
| 780 | 3-(Diethylamino)-2,2-dimethylpropyl 4-aminobenzoate and its salts <br> 3-(Diethylamino)-2,2-dimethylpropyl 4-aminobenzoate hydrochloride | - | SQE |
| 781 | N, N-diethyl-7-ethyl-4,6,6a, 7,8,9-hexahydroindolo [4,3-fg] quinoline-9-carboxamide and salts thereof | - | SQE |
| 782 | 3-[2-(diethylamino)ethyl]-1H-indol-4-ol and its salts N,N-Diethyl-4-hydroxytryptamine | - | SQE |

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|  | Chemical Material/Control Subjects | Threshold <br> (ppm) | Laws and Regulations |
| :---: | :--- | :---: | :--- |
| 783 | N, N-diethyl-7-methyl-4-propionyl-4,6,6a, 7,8,9-hexahydroindolo <br> [4,3-fg] quinoline-9-carboxamide and its salts | - | SQE |
| 784 | N,N-diethyl-7-methyl-7-methyl-4-pentanoyl-4,6,6a,7,8,9- <br> hexahydroindolo[4,3-fg]quinoline-9-carboxamide |  |  |
| 785 | N,N-diethyl-2-(5-methoxy-1H-indol-3-yl)ethanamine and its <br> salts <br> N,N-diethyl-2-(5-methoxy-1H-indol-3- <br> yl)ethanamine;hydrochloride <br> N,N-Diethyl-5-methoxytryptamine | - | SQE |
| 786 | $1-(C y c l o b u t y l m e t h y l)-N-(2-p h e n y l p r o p a n-2-y l)-1 H-i n d a z o l e-3-~$ <br> carboxamide | - | Sen |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 807 | (2,4-dimethylazetidin-1-yl)-(7-methyl-4,6,6a,7,8,9-) Hexahydroindolo[4,3-fg]quinolin-9-yl)methanone |  |  |
| 808 | 2-[(Dimethylamino) methyl] -1- (3-hydroxyphenyl) cyclohexanol and its salts | - | SQE |
| 809 | 1- (3,4-Dimethylphenyl)-2- (ethylamino) butan-1-one and its salts | - | SQE |
| 810 | 1- (3,4-Dimethylphenyl)-2- (ethylamino) pentan-1-one and its salts | - | SQE |
| 811 | 1- (3,4-Dimethylphenyl) -2- (pyrrolidin-1-yl) pentan-1-one and its salts | - | SQE |
| 812 | 2-[2,5-dimethoxy-4-(trifluoromethyl)phenyl]ethanamine and its salts <br> 2-[2,5-dimethoxy-4- <br> (trifluoromethyl)phenyl]ethanamine;hydrochloride | - | SQE |
| 813 | 2-(2,5-dimethoxy-4-nitrophenyl)ethanamine and its salts 2-(2,5-dimethoxy-4-nitrophenyl)ethanamine;hydrochloride | - | SQE |
| 814 | 1-(2,5-dimethoxy-4-nitrophenyl)propan-2-amine and its salts (2R)-1-(2,5-dimethoxy-4-nitrophenyl)propan-2-amine and its salts | - | SQE |
| 815 | 2-(2,5-dimethoxyphenyl)ethanamine and its salts <br> 2-(2,5-dimethoxyphenyl)ethanamine;hydrochloride | - | SQE |
| 816 | 1-(3,4-dimethoxyphenyl)-2-(ethylamino)pentan-1-one and its salts | - | SQE |
| 817 | 1-(3,4-dimethoxyphenyl)-2-pyrrolidin-1-ylhexan-1-one and its salts | - | SQE |
| 818 | 1-(3,4-dimethoxyphenyl)-2-pyrrolidin-1-ylpentan-1-one and its salts 1-(3,4-dimethoxyphenyl)-2-pyrrolidin-1-ylpentan-1one;hydrochloride | - | SQE |
| 819 | 1-(3,4-dimethoxyphenyl)-2-(methylamino)propan-1-one and its salts <br> 1-(3,4-dimethoxyphenyl)-2-(methylamino)propan-1one;hydrochloride | - | SQE |
| 820 | 2-(2,5-dimethoxyphenyl)-N-[(2methoxyphenyl)methyl]ethanamine and its salts | - | SQE |
| 821 | 2-(2,5-dimethoxy-4-propylphenyl)ethanamine and its salts 2-(2,5-dimethoxy-4-propylphenyl)ethanamine;hydrochloride | - | SQE |
| 822 | 1-(3,5-diMethoxy-4-propoxyphenyl)propan-2-aMine and its salts | - | SQE |
| 823 | 2-(2,5-dimethoxy-4-methylphenyl)ethanamine and its salts 2-(2,5-dimethoxy-4-methylphenyl)ethanamine;hydrochloride | - | SQE |
| 824 | 2-(2,5-dimethoxy-4-methylphenyl) -2-methoxyethaneamine and its salts | - | SQE |
| 825 | 6a,7,8,10a-Tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-yl=acetate |  |  |
| 826 | 6a,7,10,10a-Tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-yl=acetate |  |  |
| 827 | (1-((Tetrahydro-2H-pyran-4-yl)methyl)-1H-indol-3-yl)(2,2,3,3tetramethylcyclopropyl)methanone and its salts | - | SQE |
| 828 | (2,2,3,3-tetramethylcyclopropyl)-[1-(4,4,4-trifluorobutyl)indol-3yl]methanone and its salts | - | SQE |

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| :---: | :---: | :---: | :---: |
| 829 | 2-(2,4,5-trichloro-3,6-dimethoxyphenyl)ethanamine and its salts | - | SQE |
| 830 | 1-(2,4,6-trimethoxyphenyl)propan-2-amine and its salts 1-(2,4,6-trimethoxyphenyl)propan-2-amine;hydrochloride | - | SQE |
| 831 | 2- (Naphthalen-2-yl) -2- (piperidin-2-yl) acetic acid ethyl ester and its salts | - | SQE |
| 832 | ```methyl (2R)-2-naphthalen-2-yl-2-[(2R)-piperidin-2-yl]acetate and its salts methyl (2R)-2-naphthalen-2-yl-2-[(2R)-piperidin-2- yl]acetate;hydrochloride``` | - | SQE |
| 833 | 1-naphthalen-2-yl-2-pyrrolidin-1-ylpentan-1-one and its salts | - | SQE |
| 834 | naphthalen-1-yl 1-[(4-fluorophenyl)methyl]indole-3-carboxylate and its salts | - | SQE |
| 835 | Naphthalene-1-yl = 1- (5-fluoropentyl) -1H-indazole-3carboxylate and its salts | - | SQE |
| 836 | 1-(5-fluoropentyl)-N-naphthalen-1-ylindole-3-carboxamide and its salts | - | SQE |
| 837 | Naphthalene-1-yl = 1- (5-fluoropentyl) -1H-indole-3-carboxylate and its salts | - | SQE |
| 838 | naphthalen-1-yl-(1-pentylindazol-3-yl)methanone and its salts | - | SQE |
| 839 | $\mathrm{N}-1-\mathrm{Naphthalenyl}-1-$ pentyl-1H-indazole-3-carboxamide and its salts | - | SQE |
| 840 | Naphthalene-1-yl = 1-pentyl-1H-indazole-3-carboxylate and its salts | - | SQE |
| 841 | N -(Naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide and its salts | - | SQE |
| 842 | naphthalen-1-yl-(4-pentoxynaphthalen-1-yl)methanone and its salts | - | SQE |
| 843 | Naphthalen-1-yl (9-pentyl-9H-carbazol-3-yl) methanone and its salts | - | SQE |
| 844 | Naphthalen-1-yl(1-pentyl-1H-pyrrol-3-yl)methanone and its salts | - | SQE |
| 845 | Naphthalen-1-yl(1-pentyl-5-phenyl-1H-pyrrol-3-yl)methanone and its salts | - | SQE |
| 846 | N- (Naphthalen-1-yl) -1-pentyl-N- (1-pentyl-1H-indole-3carbonyl) -1 H -indole-3-carboxamide and its salts | - | SQE |
| 847 | 2-((Bis(4-fluorophenyl)methyl)sulfinyl)acetamide and its salts | - | SQE |
| 848 | 2-[Bis(4-fluorophenyl)methylsulfinyl]-N-methylacetamide and its salts | - | SQE |
| 849 | 4-Hydroxy-N-methyl-N-isopropyltryptamine and its salts | - | SQE |
| 850 | 4-Hydroxy-N,N-diisopropyltryptamine and its salts 4-Hydroxy-n,n-di(iso)propyltryptamine hydrochloride | - | SQE |
| 851 | N -\{1-[2-Hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl\}-Nphenylpropanamide | - | SQE |
| 852 | 3-(4-(1,1-Dimethylheptyl)-2-hydroxyphenyl)cyclohexanol and its salts <br> Rel-5-(1,1-Dimethylheptyl)-2-[(1R,3S)-3- <br> hydroxycyclohexyl]phenol and its salts | - | SQE |
| 853 | 3-(1-piperidin-1-ylcyclohexyl)phenol and its salts <br> 3-(1-piperidin-1-ylcyclohexyl)phenol;hydrochloride | - | SQE |
| 854 | 2- (Pyrrolidin-1-yl) -1- (thiophen-2-yl) butan-1-one and its salts | - | SQE |

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| :---: | :---: | :---: | :---: |
| 855 | 2-(Pyrrolidin-1-yl)-1-(thiophen-2-yl)pentan-1-one and its salts | - | SQE |
| 856 | 2- (Pyrrolidin-1-yl) -1- (5,6,7,8-tetrahydronaphthalen-2-yl) pentan-1-one and its salts | - | SQE |
| 857 | Propan-2-yl 2-phenyl-2-piperidin-2-ylacetate and its salts Isopropylphenidate (hydrochloride) <br> 2-yl (2R)-2-phenyl-2-[(2R)-piperidin-2-yl]acetate;hydrochloride | - | SQE |
| 858 | 1-phenyl-2-piperidin-1-ylbutan-1-one and its salts 1-phenyl-2-piperidin-1-ylbutan-1-one;hydrochloride | - | SQE |
| 859 | 1-(oxan-4-ylmethyl)-N-(2-phenylpropan-2-yl)indazole-3carboxamide and its salts | - | SQE |
| 860 | 1-(1-phenylpentan-2-yl)pyrrolidine and its salts 1-(1-phenylpentan-2-yl)pyrrolidine;hydrochloride | - | SQE |
| 861 | N - (1-phenethylpiperidin-4-yl) -Nphenylcyclopentanecarboxamide and its salts | - | SQE |
| 862 | N - (1-phenethylpiperidin-4-yl) -Nphenylcyclopentanecarboxamide and its salts | - | SQE |
| 863 | 4-Butanoyl-N, N-diethyl-7-methyl-4,6,6a, 7,8,9-hexahydroindro [4,3-fg] quinoline-9-carboxamide and its salts | - | SQE |
| 864 | 2-(butylamino) -1- (4-chlorophenyl) propan-1-one and salts thereof | - | SQE |
| 865 | 2-(Butylamino)-1-(3,4-methylenedioxyphenyl)butan-1-one |  |  |
| 866 | 2- (butylamino) -1- (3,4-methylenedioxyphenyl) pentan-1-one and salts thereof | - | SQE |
| 867 | 1-Butyl-N- (2-phenylpropan-2-yl) -1H-indole-3-carboxamide and its salts | - | SQE |
| 868 | N-\{1-[2-(Furan-2-yl)ethyl]piperidin-4-yl\}-Nphenylpropanamide | - | SQE |
| 869 | 1-(4-Fluorophenyl) -2- (isopropylamino) pentan-1-one and its salts | - | SQE |
| 870 | 1- [1- (3-fluorophenyl) cyclohexyl] piperidine and its salts | - | SQE |
| 871 | 1-(4-fluorophenyl) piperazin and its salts | - | SQE |
| 872 | methyl 2-(4-fluorophenyl)-2-(piperidin-2-yl)acetate and its salts | - | SQE |
| 873 | 1- (4-Fluorophenyl) -2- (piperidin-1-yl) pentan-1-one and its salts | - | SQE |
| 874 | N -(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2carboxamide |  |  |
| 875 | 1- (2-Fluorophenyl) Propane-2-amine and its salts | - | SQE |
| 876 | 1- (3-Fluorophenyl) Propane-2-amine and its salts | - | SQE |
| 877 | [5-(2-fluorophenyl)-1-pentylpyrrol-3-yl]-naphthalen-1ylmethanone and its salts | - | SQE |
| 878 | [5-(3-fluorophenyl)-1-pentylpyrrol-3-yl]-naphthalen-1ylmethanone and its salts | - | SQE |
| 879 | 2- (2-Fluorophenyl) -2- (methylamino) cyclohexanone and its salts | - | SQE |
| 880 | 1-(2-fluorophenyl)-N-methylpropan-2-amine and its salts | - | SQE |
| 881 | 1-(3-fluorophenyl)-N-methylpropan-2-amine and its salts 1-(3-fluorophenyl)-N-methylpropan-2-amine;hydrochloride | - | SQE |
| 882 | 1-(4-fluorophenyl)-N-methylpropan-2-amine and its salts 1-(4-fluorophenyl)-N-methylpropan-2-amine;hydrochloride | - | SQE |
| 883 | 2- (2-Fluorophenyl) -3-methylmorpholine and its salts | - | SQE |

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| :---: | :---: | :---: | :---: |
| 884 | 2－（3－fluorophenyl）－3－methylmorpholine and its salts 2－（3－fluorophenyl）－3－methylmorpholine；hydrochloride | － | SQE |
| 885 | 2－（4－Fluorophenyl）－3－methylmorpholine and its salts | － | SQE |
| 886 | ［1－（4－Fluorobenzyl）－1H－indol－3－yl］（2，2，3，3－ tetramethylcyclopropyl）methanone and its salts | － | SQE |
| 887 | ［1－（4－Fluorobenzyl）－1H－indol－3－yl］（naphthalen－1－yl） methanone and its salts | － | SQE |
| 888 | 1－（4－Fluorobenzyl）－ N －（naphthalen－1－yl）－1H－indole－3－ carboxamide and its salts | － | SQE |
| 889 | N－（2－fluorobenzyl）－2－（4－iodo－2，5－dimethoxyphenyl） ethaneamine and its salts | － | SQE |
| 890 | （1－（5－Fluoropentyl）－1H－indazol－3－yl）（naphthalen－1－yl）methanone and its salts | － | SQE |
| 891 | ［1－（5－Fluoropentyl）－1H－indol－3－yl］（pyridin－3－yl）methanone and its salts | － | SQE |
| 892 | 1－（5－Fluoropentyl）－3－（2－iodobenzoyl）indole and its salts | － | SQE |
| 893 | N －（1－Naphthyl）－1－（5－fluoropentyl）－1H－indazole－3－carboxamide and its salts | － | SQE |
| 894 | N －Phenyl－1－（5－fluoropentyl）－1H－indole－3－carboxamide and its salts | － | SQE |
| 895 | 1－（5－fluoropentyl）－N－（2－phenylpropan－2－yl）indazole－3－ carboxamide and its salts | － | SQE |
| 896 | 1－（5－fluoropentyl）－N－（2－phenylpropan－2－yl）indole－3－ carboxamide and its salts | － | SQE |
| 897 | 5－（5－Fluoropentyl）－2－（2－phenylpropan－2－yl）－2，5－dihydro－1H－ pyrido［4，3－b］indol－1－one and its salts | － | SQE |
| 898 | 1－（5－Fluoropentyl）－N－（2－phenylpropan－2－yl）－1H－pyrrolo［2，3－b］ pyridine－3－carboxamide and its salts | － | SQE |
| 899 | ［1－（5－Fluoropentyl）－1H－benzo［d］imidazol－2－yl］（naphthalen－1－ yl）methanone and its salts | － | SQE |
| 900 | 1－（4－fluoro－3－methylphenyl）－2－（pyrrolidin－1－yl）pentan－1－one |  |  |
| 901 | 2－（4－Bromo－2，5－dimethoxyphenyl）－N－（2－fluorobenzyl） ethaneamine and its salts | － | SQE |
| 902 | 2－［（4－Bromo－2，5－dimethoxyphenethylamino）methyl］phenol and its salts | － | SQE |
| 903 | 2－（8－bromo－2，3，6，7－tetrahydrobenzo［1，2－b：4，5－b＇］difuran－4－ yl ）ethanamine and its salts | － | SQE |
| 904 | 1－（8－Bromobenzo［1，2－b：4，5－b＇］difuran－4－yl）propane－2－amine and its salts | － | SQE |
| 905 | $\begin{aligned} & 6 \mathrm{a}, ~ 7, ~ 8, ~ 9, ~ 10, ~ 10 \mathrm{a} \text {-ヘキサヒドロ-6, 6, 9-トリメ } \\ & \text { チル-3-ペンチル-6 H-ジベンゾ [b, d] ピラン-1-イル = アセ } \\ & \text { テート } \end{aligned}$ |  |  |
| 906 | 6a，7，8，9，10，10a－Hexahydro－6，6，9－trimethyl－3－pentyl－6H－ dibenzo［b，d］pyran－1－ol | － | SQE |
| 907 | 2－Benzylamino－1－（3，4－methylenedioxyphenyl）propan－1－one and its salts | － | SQE |
| 908 | 4－Benzyl piperidine and its salts | － | SQE |
| 909 | N －Benzyl－1－（5－fluoropentyl）－1H－indole－3－carboxamide and its salts | － | SQE |

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| :---: | :---: | :---: | :---: |
| 910 | $\mathrm{N}-\mathrm{Benzyl}-1-\mathrm{pentyl}-1 \mathrm{H}$-indole-3-carboxamide and its salts | - | SQE |
| 911 | 1-Benzyl-4-methylpiperazine and its salts | - | SQE |
| 912 | 1-[1-(Benzo[b]thiophen-2-yl)cyclohexyl]piperidine |  | SQE |
| 913 | 1-(1-benzofuran-2-yl)-N-ethylpropan-2-amine and its salts | - | SQE |
| 914 | 1- (benzofuran-4-yl) -N-ethylpropan-2-amine and salts thereof | - | SQE |
| 915 | 1-(1-benzofuran-5-yl)-N-ethylpropan-2-amine and its salts | - | SQE |
| 916 | 1- (benzofuran-6-yl) - N -ethylpropan-2-amine and salts thereof | - | SQE |
| 917 | 1-(1-benzofuran-2-yl)propan-2-amine and its salts 1-(1-benzofuran-2-yl)propan-2-amine;hydrochloride | - | SQE |
| 918 | 1-(1-benzofuran-5-yl)propan-2-amine and its salts 1-(1-benzofuran-5-yl)propan-2-amine;hydrochloride | - | SQE |
| 919 | 1-(1-benzofuran-6-yl)propan-2-amine and its salts 1-(1-benzofuran-6-yl)propan-2-amine;hydrochloride | - | SQE |
| 920 | 1-(1-benzofuran-2-yl)-N-methylpropan-2-amine and its salts 1-(1-benzofuran-2-yl)-N-methylpropan-2-amine;hydrochloride | - | SQE |
| 921 | 1-(1-benzofuran-5-yl)-N-methylpropan-2-amine and its salts | - | SQE |
| 922 | 1-Pentyl-N- (quinolin-8-yl) -1H-indazole-3-carboxamide and its salts | - | SQE |
| 923 | 1-pentyl-N-(2-phenylpropan-2-yl)-1h-indazole-3-carboxamide and its salts | - | SQE |
| 924 | 1-pentyl-N-(2-phenylpropan-2-yl)-1h-indole-3-carboxamide and its salts | - | SQE |
| 925 | [1-(1-methylazepan-3-yl)indol-3-yl]-naphthalen-1-ylmethanone and its salts | - | SQE |
| 926 | 2- (Methylamino) -1- (3,4-dimethylphenyl) propan-1-one and its salts | - | SQE |
| 927 | 2-(Methylamino)-1-(thiophen-2-yl)propan-1-one |  | SQE |
| 928 | 2-(Methylamino)-2-phenylcyclohexanone and its salts | - | SQE |
| 929 | 2-((Methylamino)methyl)-3,4-dihydronaphthalen-1(2H)-one and its salts | - | SQE |
| 930 | N -Methylindane-2-amine and its salts | - | SQE |
| 931 | methyl (E)-2-[(2S,3S,7aS,12bS)-3-ethyl-7a-hydroxy-8-methoxy-2,3,4,6,7,12b-hexahydro-1H-indolo[2,3-a]quinolizin-2-yl]-3-methoxyprop-2-enoate and its salts | - | SQE |
| 932 | methyl (E)-2-[(2S,3S,12bS)-3-ethyl-8-methoxy-1,2,3,4,6,7,12,12b-octahydroindolo[2,3-a]quinolizin-2-yl]-3-methoxyprop-2-enoate and its salts | - | SQE |
| 933 | Methyl 2-((1-(cyclohexylmethyl)-1H-indazole-3-carbonyl)amino)-3,3-dimethylbutanoate and its salts | - | SQE |
| 934 | Methyl 2-((1-(cyclohexylmethyl)-1H-indazole-3-carbonyl)amino)-3-methylbutanoate and its salts | - | SQE |
| 935 | Methyl (2S)-2-[[1-(cyclohexylmethyl)indole-3-carbonyl]amino]-3-methylbutanoate and its salts | - | SQE |
| 936 | Methyl = 3- (3,4-dichlorophenyl) -8-methyl-8-azabicyclo [3.2.1] octane-2-carboxylate and its salts | - | SQE |
| 937 | N-Methyl-1-(2-naphthyl)propan-2-amin and its salts | - | SQE |
| 938 | (4-Methylpiperazin-1-yl) (1-pentyl-1H-indol-3-yl) methanone and its salts | - | SQE |


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| 939 | \{1-[(1-Methylpiperidin-2-yl) methyl] -1H-indole-3-yl\} (naphthalene-1-yl) methanone and its salts | - | SQE |
| 940 | 1-[1-(3-methylphenyl)cyclohexyl]pyrrolidine |  |  |
| 941 | methyl (2R)-2-(4-methylphenyl)-2-[(2S)-piperidin-2-yl]acetate and its salts | - | SQE |
| 942 | 1- (4-Methylphenyl) Propane-2-amine and its salts | - | SQE |
| 943 | Methyl (1-phenylpropan-2-yl) carbamic acid 1,1-dimethylethyl and its salts | - | SQE |
| 944 | 2-(2-Methylphenyl)-1-(1-pentyl-1H-indol-3-yl)ethanone and its salts | - | SQE |
| 945 | 1-[(3-Methylphenyl) methyl] piperazine and its salts | - | SQE |
| 946 | methyl (2S)-2-[[1-[(4-fluorophenyl)methyl]indazole-3-carbonyl]amino]-3,3-dimethylbutanoate and its salts | - | SQE |
| 947 | Methyl 2- [1- (4-fluorobenzyl) -1H-indole-3-carboxamido] -3,3dimethylbutanoate and its salts | - | SQE |
| 948 | Methyl 2-[1-(4-fluorobenzyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate | - | SQE |
| 949 | Methyl 2-[1-(4-fluorobenzyl)-1H-indole-3-carboxamido]-3methylbutanoate | - | SQE |
| 950 | Methyl 2- [1- (5-fluoropentyl) -1H-indole-3-carboxamide] -3phenylpropanoate and salts thereof | - | SQE |
| 951 | methyl (2S)-2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3methylbutanoate and its salts | - | SQE |
| 952 | Methyl 2-[1-(5-fluoropentyl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamido]-3,3-dimethylbutanoate | - | SQE |
| 953 | N-methyl-1-(5-methylthiophen-2-yl)propan-2-amine |  |  |
| 954 | 3-methyl-2-(4-methylphenyl)morpholine and its salts | - | SQE |
| 955 | Methyl = 3-Methyl-2- [1- (Penta-4-en-1-yl) -1H-Indole-3-carboxamide] Butanoate and its salts | - | SQE |
| 956 | Methyl 3-methyl-2-(((1-pentyl-1H-indazol-3$\mathrm{yl})$ carbonyl)amino)butanoate and its salts | - | SQE |
| 957 | Propylamine, N,1-dimethyl-3-(3,4-methylenedioxyphenyl)- and its salts <br> N-Methyl-3,4-methylenedioxyphenyl-3-butanamine hydrochloride | - | SQE |
| 958 | 5,6-Methylenedioxyindane-2-amine and its salts | - | SQE |
| 959 | 1-(3,4-Methylenedioxyphenyl)-2-butanamine and its salts | - | SQE |
| 960 | N-[2-(1,3-benzodioxol-5-yl)-1-methylethyl]-N-methyl-carbamicacid,1,1-dimethylethylester and its salts | - | SQE |
| 961 | 1-(3,4-Methylenedioxybenzyl) piperazine and its salts | - | SQE |
| 962 | 1-(5-methoxy-1H-indol-3-yl)propan-2-amine and its salts 1-(5-methoxy-1H-indol-3-yl)propan-2-amine;hydrochloride | - | SQE |
| 963 | (Z) -N- [3- (2-Methoxyethyl) -4,5-dimethylthiazole-2 (3H) iriden] -2,2,3,3-tetramethylcyclopropanecarboxamide and its salts | - | SQE |
| 964 | 5-Methoxy-N, N-dipropyltryptamine and its salts | - | SQE |
| 965 | 1-methoxy-3,3-dimethyl-1-oxobutane-2-yl = 1(cyclohexylmethyl) -1 H -indazole-3-carboxylate and its salts | - | SQE |
| 966 | 5-methoxy-N, N-dimethyltryptamine and its salts | - | SQE |

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| 967 | 1- [1- (4-Methoxyphenyl) cyclohexyl] piperidine and its salts | - | SQE |
| 968 | 4- [1- (3-Methoxyphenyl) cyclohexyl] morpholine and its salts | - | SQE |
| 969 | 1-(4-Methoxyphenyl) piperazin and its salts | - | SQE |
| 970 | 1-[1-(2-methoxyphenyl)-2-phenylethyl]piperidine and its salts | - | SQE |
| 971 | N - (4-Methoxyphenyl) - N - (1-phenethylpiperidin-4-yl) butanamide and its salts | - | SQE |
| 972 | 2-(3-Methoxyphenyl)-2-(propylamino)cyclohexanone | - | SQE |
| 973 | 2-(2-Methoxyphenyl)-1-(1-pentyl-1H-indol-3-yl)ethanone and its salts | - | SQE |
| 974 | (2-Methoxyphenyl)(1-pentyl-1H-indole-3-yl)methanone and its salts | - | SQE |
| 975 | (4-Methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone and its salts | - | SQE |
| 976 | 2-(3-Methoxyphenyl)-2-(methylamino) cyclohexanone and its salts | - | SQE |
| 977 | 2-(2-methoxyphenyl)-1-[1-[(1-methylpiperidin-2-yl)methyl]indol-$3-y l]$ ethanone and its salts | - | SQE |
| 978 | 3-[2-[(2-methoxyphenyl)methylamino]ethyl]-1H-quinazoline- <br> 2,4-dione and its salts | - | SQE |
| 979 | N - (2-methoxybenzyl) -2- (2,5-dimethoxy-4-methylphenyl) ethaneamine and its salts | - | SQE |
| 980 | N - (2-Methoxybenzyl) -N-methyl-1- (4-methylphenyl) propane-2-amine and its salts | - | SQE |
| 981 | 3-Methoxy-2- (methylamino) -1- (4-methylphenyl) propan-1-one and its salts | - | SQE |
| 982 | 5-methoxy-2-methyl-n,n-dimethyltryptamine and its salts | - | SQE |
| 983 | 1- (2-Methoxy-4,5-methylenedioxyphenyl) propan-2-amine and its salts | - | SQE |
| 984 | 1- (2-Methoxy-4,5-methylenedioxyphenyl) -2- (methylamino) propan-1-one and its salts | - | SQE |
| 985 | 1-(2-(4-Morpholinyl)ethyl)-3-(1-naphthoyl) indole and its salts | - | SQE |
| 986 | 5-Iodo-2-aminoindane and its salts 5-lodo-2-aminoindane hydrochloride | - | SQE |
| 987 | 1- (4-lodine-2,5-dimethoxyphenyl) propan-2-amine and its salts | - | SQE |
| 988 | 2- (4-iodo-2,5-dimethoxyphenyl) - N - (3,4methylenedioxybenzyl) ethanamine and its salts | - | SQE |
| 989 | 2-[(4-Iodine-2,5- <br> Dimethoxyphenethylamino) methyl] phenol and its salts | - | SQE |
| 990 | (2-lodine-5-nitrophenyl) \{1-[(1-methylpiperidin-2-yl) methyl] 1 H -indole-3-yl\} metanone and its salts | - | SQE |
| 991 | (2-lodophenyl)(1-pentyl-1H-indol-3-yl)methanone and its salts | - | SQE |
| 992 | (2-iodophenyl)-[1-(1-methylazepan-3-yl)indol-3-yl]methanone and its salts | - | SQE |
| 993 | (2-lodophenyl) \{1-[(1-methylpiperidin-2-yl) methyl] -1H-indole-$3-y \mid\}$ metanone and its salts | - | SQE |
| 994 | (2-Iodophenyl)(1-pentyl-1H-indol-3-yl)methanone and its salts | - | SQE |
| 995 | (2-Iodophenyl)[1-(1-methylazepan-3-yl)-1H-indole 1 H -indole-$3-y l] m e t h a n o n e ~ a n d ~ i t s ~ s a l t s ~$ | - | SQE |

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| 996 | (2-lodophenyl)\{1-[(1-methylpiperidin-2-yl)methyl]-1H-indole methanone and its salts | - | SQE |
| 997 | (1H-Indol-3-yl) (naphthalen-1-yl) methanone | - | SQE |
| 998 | (2-Methyl-1H-indol-3-yl) (naphthalen-1-yl) methanone | - | SQE |
| 999 | 2-amino-1-phenyl-propan-1-one | - | SQE |
| 1000 | 6a,7,8,10a-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran-1-ol | - | SQE |
| 1001 | 6a,7,10,10a-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran-1ol | - | SQE |
| 1002 | A substance containing any of those listed above. However, plants other than Salvia divinoram(it is limited to the one that can be used for the human body immediately) and Mitragyna speciosa(including hybrids with other species belonging to Mitragyna, and be limited to the one that can be used for the human body immediately) are excluded. | - | SQE |
| 1003 | 5-tert-butyl-2,4,6- trinitro-m-xylene (musk xylene) | - | REACH Annex XIV |
| 1004 | 4,4'- Diaminodiphenylmethane (MDA) | - | REACH Annex XIV |
| 1005 | Bis (2-ethyl(hexyl)phthalate) (DEHP) | - | REACH Annex XIV |
| 1006 | Benzyl butyl phthalate (BBP) | - | REACH Annex XIV |
| 1007 | Dibutyl phthalate (DBP) | - | REACH Annex XIV |
| 1008 | hexabromocyclododecane (HBCDD) [alpha-hexabromocyclododecane, beta-hexabromocyclododecane, gamma- hexabromocyclododecane] | - | REACH Annex XIV |
| 1009 | Diisobutyl phthalate (DIBP) | - | REACH Annex XIV |
| 1010 | Diarsenic trioxide | - | REACH Annex XIV |
| 1011 | Diarsenic pentaoxide | - | REACH Annex XIV |
| 1012 | Lead chromate | - | REACH Annex XIV |
| 1013 | Lead sulfochromate yellow | - | REACH Annex XIV |
| 1014 | Lead chromate molybdate sulphate red | - | REACH Annex XIV |
| 1015 | Tris (2-chloroethyl) phosphate (TCEP) | - | REACH Annex XIV |
| 1016 | 2,4-Dinitrotoluene (2,4-DNT) | - | REACH Annex XIV |
| 1017 | Trichloroethylene | - | REACH Annex XIV |
| 1018 | Chromium trioxide | - | REACH Annex XIV |
| 1019 | Acids generated from chromium trioxide and their oligomers Group containing: <br> Chromic acid Dichromic acid Oligomers of chromic acid and dichromic acid | - | REACH Annex XIV |
| 1020 | Sodium dichromate | - | REACH Annex XIV |
| 1021 | Potassium dichromate | - | REACH Annex XIV |
| 1022 | Ammonium dichromate | - | REACH Annex XIV |
| 1023 | Potassium chromate | - | REACH Annex XIV |
| 1024 | Sodium chromate | - | REACH Annex XIV |
| 1025 | Formaldehyde, oligomeric reaction products with aniline | - | REACH Annex XIV |
| 1026 | Arsenic acid | - | REACH Annex XIV |
| 1027 | Bis(2-methoxyethyl) ether (diglyme) | - | REACH Annex XIV |
| 1028 | 1,2-dichloroethane (EDC) | - | REACH Annex XIV |

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| 1029 | 2,2' -dichloro-4,4' -methylenedianiline (MOCA) | - | REACH Annex XIV |
| 1030 | Dichromium tris(chromate) | - | REACH Annex XIV |
| 1031 | Strontium chromate | - | REACH Annex XIV |
| 1032 | Potassium hydroxyoctaoxodizincatedichromate | - | REACH Annex XIV |
| 1033 | Pentazinc chromate octahydroxide | - | REACH Annex XIV |
| 1034 | 1-Bromopropane (n-propyl bromide) | - | REACH Annex XIV |
| 1035 | Diisopentylphthalate | - | REACH Annex XIV |
| 1036 | 1,2-Benzenedicarboxylic acid, di-C6-8-branched alkyl esters, C7 rich | - | REACH Annex XIV |
| 1037 | 1,2-Benzenedicarboxylic acid, di-C7-11-branched and linear alkyl esters | - | REACH Annex XIV |
| 1038 | 1,2-Benzenedicarboxylic acid, dipentylester, branched and linear | - | REACH Annex XIV |
| 1039 | Bis(2-methoxyethyl) phthalate | - | REACH Annex XIV |
| 1040 | Dipentyl phthalate (DPP) | - | REACH Annex XIV |
| 1041 | N-pentyl-isopentylphthalate | - | REACH Annex XIV |
| 1042 | Anthracene oil | - | REACH Annex XIV |
| 1043 | Pitch, coal tar, high temp. | - | REACH Annex XIV |
| 1044 | 4-(1,1,3,3-Tetramethylbutyl) phenol, ethoxylated (covering well-defined substances and UVCB substances, polymers and homologues) | - | REACH Annex XIV |
| 1045 | 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 welldefined substances, polymers and homologues, which include any of the individual isomers and/or combinations thereof) | - | REACH Annex XIV |
| 1046 | 1,2-Benzenedicarboxylic acid, dihexyl ester, branched and linear | - | REACH Annex XIV |
| 1047 | Dihexyl phthalate | - | REACH Annex XIV |
| 1048 | 1,2-benzenedicarboxylic acid, di-C6-10-alkyl esters or mixed decyl and hexyl and octyl diesters(with $\geq 0.3 \%$ of dihexyl phthalate (EC No. 201-559-5)) | - | REACH Annex XIV |
| 1049 | Trixylyl phosphate | - | REACH Annex XIV |
| 1050 | Sodium perborate, perboric acid, sodium salt | - | REACH Annex XIV |
| 1051 | Sodium peroxometaborate | - | REACH Annex XIV |
| 1052 | 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-1,3dioxane [2] (covering any of the individual stereoisomers of [1] and [2] or any combination thereof) | - | REACH Annex XIV |
| 1053 | 2-(2H-benzotriazol-2-yl)-4,6-ditertpentylphenol (UV-328) | - | REACH Annex XIV |
| 1054 | 2,4-di-tert-butyl-6-(5-chlorobenzotriazol-2-yl)phenol(UV-327) | - | REACH Annex XIV |
| 1055 | 2-(2H-benzotriazol-2-yl)-4-(tert-butyl)-6-(sec-butyl)phenol | - | REACH Annex XIV |
| 1056 | 2-benzotriazol-2-yl-4,6-di-tert-butylphenol(UV-320) | - | REACH Annex XIV |
| 1057 | Tetraethyllead | - | REACH Annex XIV |
| 1058 | 4,4'-bis(dimethylamino)-4"-(methylamino)trityl alcohol (with $\geqq 0.1 \%$ of Michler's ketone (EC No. 202-027-5) or Michler's base (EC No. 202-959-2)) | - | REACH Annex XIV |
| 1059 | Reaction products of 1,3,4-thiadiazolidine-2,5-dithione, formaldehyde and 4-heptylphenol, branched and linear (RP-HP) <br> [with $\geqq 0.1 \% \mathrm{w} / \mathrm{w} 4$-heptylphenol, branched and linear] | - | REACH Annex XIV |
| 1060 | 2-ethylhexyl 10-ethyl-4,4-dioctyl-7-oxo-8-oxa-3,5-dithia-4stannatetradecanoate (DOTE) | - | REACH Annex XIV |

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| 1061 | 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) | - | REACH Annex XIV |
| 1062 | Polychlorinated terphenyls (PCTs) | - | REACH Annex XVII |
| 1063 | Chloroethene, (Vinyl chloride) | - | REACH Annex XVII |
| 1064 | Liquid substances or mixtures which are regarded as dangerous in accordance with Directive 1999/45/EC or are fulfilling the criteria for any of the following hazard classes or categories set out in Annex I to Regulation (EC) No 1272/2008: | - | REACH Annex XVII |
| 1065 | Tris (2,3 dibromopropyl) phosphate | - | REACH Annex XVII |
| 1066 | Benzene | - | REACH Annex XVII |
| 1067 | Asbestos fibres | - | REACH Annex XVII |
| 1068 | Tris(aziridinyl)phosphinoxide | - | REACH Annex XVII |
| 1069 | Polybromobiphenyls, Polybrominatedbiphenyls (PBB) | - | REACH Annex XVII |
| 1070 | Soap bark powder (Quillaja saponaria) and its derivatives containing saponines | - | REACH Annex XVII |
| 1071 | Powder of the roots of Helleborus viridis and Helleborus niger | - | REACH Annex XVII |
| 1072 | Powder of the roots of Veratrum album and Veratrum nigrum | - | REACH Annex XVII |
| 1073 | Benzidine and/or its derivatives | - | REACH Annex XVII |
| 1074 | o-Nitrobenzaldehyde | - | REACH Annex XVII |
| 1075 | Wood powder | - | REACH Annex XVII |
| 1076 | Ammonium sulphide | - | REACH Annex XVII |
| 1077 | Ammonium hydrogen sulphide | - | REACH Annex XVII |
| 1078 | Ammonium polysulphide | - | REACH Annex XVII |
| 1079 | Volatile esters of bromoacetic acids | - | REACH Annex XVII |
| 1080 | 2-naphthylamine and its salts | - | REACH Annex XVII |
| 1081 | Benzidine and its salts | - | REACH Annex XVII |
| 1082 | 4-Nitrobiphenyl | - | REACH Annex XVII |
| 1083 | 4-Aminobiphenyl xenylamine and its salts | - | REACH Annex XVII |
| 1084 | Lead carbonates | - | REACH Annex XVII |
| 1085 | Lead sulphates | - | REACH Annex XVII |
| 1086 | Mercury compounds | - | REACH Annex XVII |
| 1087 | Mercury | - | REACH Annex XVII |
| 1088 | Arsenic compounds | - | REACH Annex XVII |
| 1089 | Organostannic compounds | - | REACH Annex XVII |
| 1090 | Di- $\mu$-oxo-di-n-butylstanniohydroxyborane / Dibutyltin hydrogen borate C8H19BO3Sn (DBB) | - | REACH Annex XVII |
| 1091 | Cadmium and its compounds | - | REACH Annex XVII |
| 1092 | Monomethyl-tetrachlorodiphenyl methane Trade name: Ugilec 141 | - | REACH Annex XVII |
| 1093 | Monomethyl-dichloro-diphenyl methane Trade name: Ugilec 121, Ugilec 21 | - | REACH Annex XVII |
| 1094 | Monomethyl-dibromo-diphenyl methane bromobenzylbromotoluene, mixture of isomers Trade name: DBBT | - | REACH Annex XVII |
| 1095 | Nickel and its compounds | - | REACH Annex XVII |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 1096 | Substances which are classified as carcinogen category 1A or 1B in Part 3 of Annex VI to Regulation (EC) No 1272/2008 and are listed in Appendix 1 or Appendix 2, respectively. | - | REACH Annex XVII |
| 1097 | Substances which are classified as germ cell mutagen category 1A or 1B in Part 3 of Annex VI to Regulation (EC) No 1272/2008 and are listed in Appendix 3 or Appendix 4, respectively. | - | REACH Annex XVII |
| 1098 | Substances which are classified as reproductive toxicant category 1A or 1B in Part 3 of Annex VI to Regulation (EC) No 1272/2008 and are listed in Appendix 5 or Appendix 6, respectively | - | REACH Annex XVII |
| 1099 | Creosote and Creosote related substances | - | REACH Annex XVII |
| 1100 | Chloroform | - | REACH Annex XVII |
| 1101 | 1,1,2-Trichloroethane | - | REACH Annex XVII |
| 1102 | 1,1,2,2-Tetrachloroethane | - | REACH Annex XVII |
| 1103 | 1,1,1,2-Tetrachloroethane | - | REACH Annex XVII |
| 1104 | Pentachloroethane | - | REACH Annex XVII |
| 1105 | 1,1-Dichloroethene | - | REACH Annex XVII |
| 1106 | Substances classified as flammable gases category 1 or 2 , flammable liquids categories 1 , 2 or 3 , flammable solids category 1 or 2 ,substances and mixtures which, in contact with water, emit flammable gases, category 1, 2 or 3 , pyrophoric liquids category 1 or pyrophoric solids category 1 , regardless of whether they appear in Part 3 of Annex VI to Regulation (EC) No 1272/2008 or not | - | REACH Annex XVII |
| 1107 | Hexachloroethane | - | REACH Annex XVII |
| 1108 | Azocolourants and Azodyes | - | REACH Annex XVII |
| 1109 | Diphenylether, octabromo derivative C 12 H 2 Br 8 O | - | REACH Annex XVII |
| 1110 | Nonylphenol C6H4(OH)C9H19 | - | REACH Annex XVII |
| 1111 | Nonylphenol ethoxylates ( C 2 H 4 O ) $\mathrm{nC15H24O}$ | - | REACH Annex XVII |
| 1112 | Chromium VI compounds | - | REACH Annex XVII |
| 1113 | Toluene | - | REACH Annex XVII |
| 1114 | Trichlorobenzene | - | REACH Annex XVII |
| 1115 | Polycyclic-aromatic hydrocarbons (PAH) | - | REACH Annex XVII |
| 1116 | The following phthalates (or other CAS and EC numbers covering the substance) | - | REACH Annex XVII |
| 1117 | Bis (2-ethylhexyl) phthalate (DEHP) | - | REACH Annex XVII |
| 1118 | Dibutyl phthalate (DBP) | - | REACH Annex XVII |
| 1119 | Benzyl butyl phthalate (BBP) | - | REACH Annex XVII |
| 1120 | Diisobutyl phthalate (DIBP) | - | REACH Annex XVII |
| 1121 | The following phthalates (or other CAS and EC numbers covering the substance) | - | REACH Annex XVII |
| 1122 | Di-isononyl phthalate (DINP) | - | REACH Annex XVII |
| 1123 | Di-isodecyl phthalate (DIDP) | - | REACH Annex XVII |
| 1124 | Di-n-octyl phthalate (DNOP) | - | REACH Annex XVII |
| 1125 | 2-(2-methoxyethoxy)ethanol (DEGME) | - | REACH Annex XVII |
| 1126 | 2-(2-butoxyethoxy)ethanol (DEGBE) | - | REACH Annex XVII |
| 1127 | Methylenediphenyl diisocyanate (MDI) including the following specific isomers | - | REACH Annex XVII |
| 1128 | 4,4'-Methylenediphenyl diisocyanate | - | REACH Annex XVII |
| 1129 | 2,4'-Methylenediphenyl diisocyanate | - | REACH Annex XVII |

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|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 1130 | 2,2'-Methylenediphenyl diisocyanate | - | REACH Annex XVII |
| 1131 | Cyclohexane | - | REACH Annex XVII |
| 1132 | Ammonium nitrate (AN) | - | REACH Annex XVII |
| 1133 | Dichloromethane | - | REACH Annex XVII |
| 1134 | Acrylamide | - | REACH Annex XVII |
| 1135 | Dimethylfumarate (DMF) | - | REACH Annex XVII |
| 1136 | Phenylmercury acetate | - | REACH Annex XVII |
| 1137 | Phenylmercury propionate | - | REACH Annex XVII |
| 1138 | Phenylmercury 2-ethylhexanoate | - | REACH Annex XVII |
| 1139 | Phenylmercury octanoate | - | REACH Annex XVII |
| 1140 | Phenylmercury neodecanoate | - | REACH Annex XVII |
| 1141 | Lead and its compounds | - | REACH Annex XVII |
| 1142 | 1,4-Dichlorobenzene | - | REACH Annex XVII |
| 1143 | Inorganic ammonium salts | - | REACH Annex XVII |
| 1144 | 4,4'-isopropylidenediphenol Bisphenol A; BPA | - | REACH Annex XVII |
| 1145 | C9-C14 linear and/or branched perfluorocarboxylic acids (C9-C14 PFCAs), their salts and C9-C14 PFCAs-related substances, perfluorononan-1-oic acid (PFNA); nonadecafluorodecanoic acid (PFDA); henicosafluoroundecanoic acid (PFUnDA); tricosafluorododecanoic acid (PFDoDA); pentacosafluorotridecanoic acid (PFTrDA); heptacosafluorotetradecanoic acid (PFTDA); including their salts and precursors | - | REACH Annex XVII |
| 1146 | Methanol | - | REACH Annex XVII |
| 1147 | Octamethylcyclotetrasiloxane (D4) | - | REACH Annex XVII |
| 1148 | Decamethylcyclopentasiloxane (D5) | - | REACH Annex XVII |
| 1149 | 1-methyl-2-pyrrolidone | - | REACH Annex XVII |
| 1150 | The following substances which are classified as carcinogenic, mutagenic or toxic for reproduction, category 1A or 1B (See group members) <br> The substances listed in column 1 of the Table in Appendix 12 | - | REACH Annex XVII |
| 1151 | ( $3,3,4,4,5,5,6,6,7,7,8,8,8$-tridecafluorooctyl) silanetriol and Any of its mono-, di- or tri-O-(alkyl) derivatives (TDFAs) | - | REACH Annex XVII |
| 1152 | Diisocyanates |  | REACH Annex XVII |
| 1153 | Substances in tattoo inks and permanent make up |  | REACH Annex XVII |
| 1154 | N,N-Dimethylformamide |  | REACH Annex XVII |
| 1155 | Formaldehyde and formaldehyde-releasing substances |  | REACH Annex XVII |
| 1156 | synthetic polymer microparticles |  | REACH Annex XVII |
| 1157 | 2,4,5-T and its salts and esters | - | PIC Convention |
| 1158 | Alachlor | - | PIC Convention |
| 1159 | Aldicarb | - | PIC Convention |
| 1160 | Aldrin | - | PIC Convention |
| 1161 | Azinphos-methyl | - | PIC Convention |
| 1162 | Binapacryl | - | PIC Convention |
| 1163 | Captafol | - | PIC Convention |
| 1164 | Carbofuran | - | PIC Convention |
| 1165 | Chlordane | - | PIC Convention |
| 1166 | Chlordimeform | - | PIC Convention |


|  | Chemical Material/Control Subjects | Threshold (ppm) | Laws and Regulations |
| :---: | :---: | :---: | :---: |
| 1167 | Chlorobenzilate | - | PIC Convention |
| 1168 | DDT | - | PIC Convention |
| 1169 | Dieldrin | - | PIC Convention |
| 1170 | Dinitro-ortho-cresol (DNOC) and its salts (such as ammonium salt, potassium salt and sodium salt) | - | PIC Convention |
| 1171 | Dinoseb and its salts and esters | - | PIC Convention |
| 1172 | EDB (1,2-dibromoethane) | - | PIC Convention |
| 1173 | Endosulfan | - | PIC Convention |
| 1174 | Ethylene dichloride | - | PIC Convention |
| 1175 | Ethylene oxide | - | PIC Convention |
| 1176 | Fluoroacetamide | - | PIC Convention |
| 1177 | HCH (mixed isomers) | - | PIC Convention |
| 1178 | Heptachlor | - | PIC Convention |
| 1179 | Hexachlorobenzene | - | PIC Convention |
| 1180 | Lindane (gamma-HCH) | - | PIC Convention |
| 1181 | Mercury compounds, including inorganic mercury compounds, alkyl mercury compounds and alkyloxyalkyl and aryl mercury compounds | - | PIC Convention |
| 1182 | Methamidophos | - | PIC Convention |
| 1183 | Monocrotophos | - | PIC Convention |
| 1184 | Parathion | - | PIC Convention |
| 1185 | Pentachlorophenol and its salts and esters | - | PIC Convention |
| 1186 | Toxaphene (Camphechlor) | - | PIC Convention |
| 1187 | Tributyl tin compounds | - | PIC Convention |
| 1188 | Trichlorfon | - | PIC Convention |
| 1189 | Dustable powder formulations containing a combination of benomyl at or above 7\%, carbofuran at or above $10 \%$ and thiram at or above 15\% | - | PIC Convention |
| 1190 | Phosphamidon | - | PIC Convention |
| 1191 | Methyl-parathion | - | PIC Convention |
| 1192 | Asbestos(Actinolite,Anthophyllite,Amosite,Crocidolite,Tremolite) | - | PIC Convention |
| 1193 | Commercial octabromodiphenyl ether (including Hexabromodiphenyl ether and Heptabromodiphenyl ether) | - | PIC Convention |
| 1194 | Commercial pentabromodiphenyl ether (including tetrabromodiphenyl ether and pentabromodiphenyl ether) | - | PIC Convention |
| 1195 | Perfluorooctane sulfonic acid, perfluorooctane sulfonates, perfluorooctane sulfonamides and perfluorooctane sulfonyls | - | PIC Convention |
| 1196 | Polybrominated Biphenyls (PBBs) | - | PIC Convention |
| 1197 | Polychlorinated Biphenyls (PCBs) | - | PIC Convention |
| 1198 | Polychlorinated Terphenyls (PCTs) | - | PIC Convention |
| 1199 | Short-chain chlorinated paraffins (SCCP) | - | PIC Convention |
| 1200 | Tetraethyl lead | - | PIC Convention |
| 1201 | Tetramethyl lead | - | PIC Convention |
| 1202 | Tributyl tin compounds | - | PIC Convention |
| 1203 | Tris(2,3 dibromopropyl)phosphate | - | PIC Convention |

("-" means there is no intentional addition (component).)
Abbreviation of the Laws:

| ISHA: | Industrial Safety and Health Act of Japan <br> ECS: |
| :--- | :--- |
| RNSM: | Act on the Regulation of Manufacture and Evaluation of Chemical Substances of Japan <br> of Japan |
| PCW: | Act on the Prohibition of Chemical Weapons and the Regulation of Specific Chemicals <br> of Japan |
| PDSC: | Poisonous and Deleterious Substances Control Act of Japan |
| POL: | Act on the Protection of the Ozone Layer Through the Control of Specified <br> Substances, etc. and Other Measures of Japan |
| NPC: | Narcotics and Psychotropic Control Act of Japan |
| PGWC: | Act on Promotion of Global Warming Countermeasures of Japan <br> Act on Securing Quality, Efficacy and Safety of Products Including Pharmaceuticals <br> and Medical Devices of Japan |
| PEPM: | Act on Preventing Environmental Pollution of Mercury of Japan |

