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HOUSE BILL NO. 193

Offered January 12, 2022

Prefiled January 10, 2022

A *BILL to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Drug Control Act; Schedule I.*

 Patron—Hodges

Committee Referral Pending

Be it enacted by the General Assembly of Virginia:**1. That § 54.1-3446 of the Code of Virginia is amended and reenacted as follows:****§ 54.1-3446. Schedule I.**

The controlled substances listed in this section are included in Schedule I:

1. Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers and salts is possible within the specific chemical designation:

1-[1-[1-(4-bromophenyl)ethyl]-4-piperidinyl]-1,3-dihydro-2H-benzimidazol-2-one (other name: Biorphine);

1-[2-methyl-4-(3-phenyl-2-propen-1-yl)-1-piperazinyl]-1-butanone (other name: 2-methyl AP-237);

1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP);

1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP);

2-[(4-methoxyphenyl)methyl]-N,N-diethyl-5-nitro-1H-benzimidazole-1-ethanamine (other name: Metonitazene);

2-methoxy-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide (other name: Methoxyacetyl fentanyl);

3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzamide (other name: U-47700);

3,4-dichloro-N-[[1-(dimethylamino)cyclohexyl]methyl]benzamide (other name: AH-7921);

Acetyl fentanyl (other name: desmethyl fentanyl);

Acetylmethadol;

Allylprodine;

Alphacetylmethadol (except levo-alpha-cetylmethadol, also known as levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM);

Alphameprodine;

Alphamethadol;

Benzethidine;

Betacetylmethadol;

Betameprodine;

Betamethadol;

Betaprodine;

Clonitazene;

Dextromoramide;

Diampromide;

Diethylthiambutene;

Difenoxin;

Dimenoxadol;

Dimepheptanol;

Dimethylthiambutene;

Dioxaphetylbutyrate;

Dipipanone;

Ethylmethylthiambutene;

Etonitazene;

Etoxidine;

Furethidine;

Hydroxypethidine;

Ketobemidone;

Levomoramide;

Levophenacetylmorphan;

Morpheridine;

INTRODUCED

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- 59 MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
60 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (other name: Cyclopropyl fentanyl);
61 N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (other name: Tetrahydrofuranyl
62 fentanyl);
63 N-[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]-N-phenylpropanamide (other name:
64 alpha-methylthiofentanyl);
65 N-[1-(1-methyl-2-phenylethyl)-4-piperidyl]-N-phenylacetamide (other name:
66 acetyl-alpha-methylfentanyl);
67 N-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyll}-N-phenylpropanamide (other name:
68 beta-hydroxythiofentanyl);
69 N-[1-(2-hydroxy-2-phenyl)ethyl-4-piperidyl]-N-phenylpropanamide (other name:
70 beta-hydroxyfentanyl);
71 N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide (other names:
72 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine, alpha-methylfentanyl);
73 N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other names: 2-fluorofentanyl,
74 ortho-fluorofentanyl);
75 N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name: 3-fluorofentanyl);
76 N-[3-methyl-1-(2-hydroxy-2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (other name:
77 beta-hydroxy-3-methylfentanyl);
78 N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (other name: 3-methylfentanyl);
79 N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyll]-N-phenylpropanamide (other name:
80 3-methylthiofentanyl);
81 *N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other names:*
82 *para-chlorofentanyl, 4-chlorofentanyl);*
83 N-(4-fluorophenyl)-2-methyl-N-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name:
84 para-fluoroisobutyl fentanyl);
85 N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyll]-butanamide (other name:
86 para-fluorobutylfentanyl);
87 N-(4-fluorophenyl)-N-1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name: para-fluorofentanyl);
88 N,N-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine (other name:
89 Isotonitazene);
90 *N,N-diethyl-2-[(4-ethoxyphenyl) methyl]-1H-benzimidazol-1-yl}-ethan-1-amine (other names: Etazene,*
91 *Desnitroetonitazene);*
92 *N,N-diethyl-2-[(4-methoxyphenyl)methyl]-1H-benzimidazole-1-ethanamine (other name:*
93 *Metodesnitazene);*
94 N-phenyl-N-[1-(2-phenylmethyl)-4-piperidinyll]-2-furancarboxamide (other name: N-benzyl Furanyl
95 norfentanyl);
96 N-phenyl-N-(4-piperidinyll)-propanamide (other name: Norfentanyl);
97 Noracymethadol;
98 Norlevorphanol;
99 Normethadone;
100 Norpipanone;
101 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyll]-2-furancarboxamide (other name: Furanyl fentanyl);
102 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyll]-2-propenamide (other name: Acryl fentanyl);
103 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyll]-butanamide (other name: butyl fentanyl);
104 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyll]-pentanamide (other name: Pentanoyl fentanyl);
105 N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyll]-propanamide (other name: thiofentanyl);
106 Phenadoxone;
107 Phenampromide;
108 Phenomorphan;
109 Phenoperidine;
110 Piritramide;
111 Proheptazine;
112 Properidine;
113 Propiram;
114 Racemoramide;
115 Tilidine;
116 Trimeperidine;
117 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyll]-1,3-benzodioxole-5-carboxamide (other name:
118 Benzodioxole fentanyl);
119 3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methylbenzamide (other name: U-49900);
120 2-(2,4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methyl acetamide (other name: U-48800);

- 121 2-(3,4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methyl acetamide (other name: U-51754);
 122 N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)-4-piperidiny]-acetamide (other name: Ocfentanil);
 123 N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidiny]-butanamide (other name:
 124 4-methoxybutyrylfentanyl);
 125 N-phenyl-2-methyl-N-[1-(2-phenylethyl)-4-piperidiny]-propanamide (other name: Isobutyryl fentanyl);
 126 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-cyclopentanecarboxamide (other name: Cyclopentyl
 127 fentanyl);
 128 N-phenyl-N-(1-methyl-4-piperidiny)-propanamide (other name: N-methyl norfentanyl);
 129 N-[2-(dimethylamino)cyclohexyl]-N-methyl-1,3-benzodioxole-5-carboxamide (other names:
 130 3,4-methylenedioxy U-47700 or 3,4-MDO-U-47700);
 131 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-2-butenamide (other name: Crotonyl fentanyl);
 132 N-phenyl-N-[4-phenyl-1-(2-phenylethyl)-4-piperidiny]-propanamide (other name: 4-phenylfentanyl);
 133 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-benzamide (other names: Phenyl fentanyl, Benzoyl
 134 fentanyl);
 135 N-[2-(dimethylamino)cyclohexyl]-N-phenylfuran-2-carboxamide (other name: Furanyl UF-17);
 136 N-[2-(dimethylamino)cyclohexyl]-N-phenylpropionamide (other name: UF-17);
 137 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-isopropyl-benzamide (other name: Isopropyl
 138 U-47700).
 139 2. Any of the following opium derivatives, their salts, isomers and salts of isomers, unless
 140 specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible
 141 within the specific chemical designation:
 142 Acetorphine;
 143 Acetyldihydrocodeine;
 144 Benzylmorphine;
 145 Codeine methylbromide;
 146 Codeine-N-Oxide;
 147 Cyprenorphine;
 148 Desomorphine;
 149 Dihydromorphine;
 150 Drotebanol;
 151 Etorphine;
 152 Heroin;
 153 Hydromorphenol;
 154 Methyldesorphine;
 155 Methyldihydromorphine;
 156 Morphine methylbromide;
 157 Morphine methylsulfonate;
 158 Morphine-N-Oxide;
 159 Myrophine;
 160 Nicocodeine;
 161 Nicomorphine;
 162 Normorphine;
 163 Pholcodine;
 164 Thebacon.
 165 3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture,
 166 or preparation, which contains any quantity of the following hallucinogenic substances, or which
 167 contains any of its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers,
 168 and salts of isomers is possible within the specific chemical designation (for purposes of this subdivision
 169 only, the term "isomer" includes the optical, position, and geometric isomers):
 170 Alpha-ethyltryptamine (some trade or other names: Monase; a-ethyl-1H-indole-3-ethanamine;
 171 3-2-aminobutyl] indole; a-ET; AET);
 172 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:
 173 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B; Nexus);
 174 3,4-methylenedioxy amphetamine;
 175 5-methoxy-3,4-methylenedioxy amphetamine;
 176 3,4,5-trimethoxy amphetamine;
 177 Alpha-methyltryptamine (other name: AMT);
 178 Bufotenine;
 179 Diethyltryptamine;
 180 Dimethyltryptamine;
 181 4-methyl-2,5-dimethoxyamphetamine;

- 182 2,5-dimethoxy-4-ethylamphetamine (DOET);
 183 4-fluoro-N-ethylamphetamine;
 184 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
 185 Ibogaine;
 186 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
 187 Lysergic acid diethylamide;
 188 Mescaline;
 189 Parahexyl (some trade or other names:
 190 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl);
 191 Peyote;
 192 N-ethyl-3-piperidyl benzilate;
 193 N-methyl-3-piperidyl benzilate;
 194 Psilocybin;
 195 Psilocyn;
 196 Salvinorin A;
 197 Tetrahydrocannabinols, except as present in (i) industrial hemp, as defined in § 3.2-4112, that is
 198 possessed by a person registered pursuant to subsection A of § 3.2-4115 or his agent; (ii) a hemp
 199 product, as defined in § 3.2-4112, containing a tetrahydrocannabinol concentration of no greater than 0.3
 200 percent that is derived from industrial hemp, as defined in § 3.2-4112, that is grown, dealt, or processed
 201 in compliance with state or federal law; (iii) marijuana; (iv) dronabinol in sesame oil and encapsulated
 202 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration; or (v)
 203 industrial hemp, as defined in § 3.2-4112, that is possessed by a person who holds a hemp producer
 204 license issued by the U.S. Department of Agriculture pursuant to 7 C.F.R. Part 990;
 205 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;
 206 2,5-DMA);
 207 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts
 208 and salts of isomers;
 209 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4
 210 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);
 211 N-hydroxy-3,4-methylenedioxyamphetamine (some other names:
 212 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);
 213 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:
 214 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);
 215 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;
 216 paramethoxyamphetamine; PMA);
 217 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,
 218 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);
 219 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy,
 220 PHP);
 221 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine,
 222 2-thienyl analog of phencyclidine, TPCP, TCP);
 223 1-1-(2-thienyl)cyclohexylpyrrolidine (other name: TCPy);
 224 3,4-methylenedioxypyrovalerone (other name: MDPV);
 225 4-methylmethcathinone (other names: mephedrone, 4-MMC);
 226 3,4-methylenedioxymethcathinone (other name: methylone);
 227 Naphthylpyrovalerone (other name: naphyrone);
 228 4-fluoromethcathinone (other names: flephedrone, 4-FMC);
 229 4-methoxymethcathinone (other names: methedrone; bk-PMMA);
 230 Ethcathinone (other name: N-ethylcathinone);
 231 3,4-methylenedioxyethcathinone (other name: ethylone);
 232 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone);
 233 N,N-dimethylcathinone (other name: metamfepramone);
 234 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);
 235 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);
 236 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);
 237 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);
 238 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);
 239 3-fluoromethcathinone (other name: 3-FMC);
 240 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);
 241 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);
 242 4-Methylethcathinone (other name: 4-MEC);
 243 4-Ethylmethcathinone (other name: 4-EMC);

- 244 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);
 245 Beta-keto-methylbenzodioxolypentanamine (other names: Pentylone, bk-MBDP);
 246 Alpha-methylamino-butyrophenone (other name: Buphedrone);
 247 Alpha-methylamino-valerophenone (other name: Pentedrone);
 248 3,4-Dimethylmethcathinone (other name: 3-,4-DMMC);
 249 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);
 250 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,
 251 25I-NBOMe, 2C-I-NBOMe);
 252 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);
 253 4-Fluoromethamphetamine (other name: 4-FMA);
 254 4-Fluoroamphetamine (other name: 4-FA);
 255 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);
 256 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);
 257 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);
 258 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);
 259 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);
 260 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
 261 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
 262 (2-aminopropyl)benzofuran (other name: APB);
 263 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
 264 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 265 2C-C-NBOMe, 25C-NBOMe, 25C);
 266 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 267 2C-B-NBOMe, 25B-NBOMe, 25B);
 268 Acetoxymethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
 269 Benocyclidine (other names: BCP, BTCP);
 270 Alpha-pyrrolidinobutyrophenone (other name: alpha-PBP);
 271 3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
 272 4-bromomethcathinone (other name: 4-BMC);
 273 4-chloromethcathinone (other name: 4-CMC);
 274 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
 275 Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
 276 Alpha-Pyrrolidinoheptiophenone (other name: PV8);
 277 5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
 278 Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
 279 Beta-keto-4-bromo-2,5-dimethoxyphenethylamine (other name: bk-2C-B);
 280 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
 281 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
 282 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
 283 4-Chloroethcathinone (other name: 4-CEC);
 284 3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
 285 1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
 286 (2-Methylaminopropyl)benzofuran (other name: MAPB);
 287 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)-1-pentanone (other names: N,N-Dimethylpentylone,
 288 Dipentylone);
 289 1-(4-methoxyphenyl)-2-(pyrrolidin-1-yl)octan-1-one (other name: 4-methoxy-PV9);
 290 3,4-tetramethylene-alpha-pyrrolidinovalerophenone (other name: TH-PVP);
 291 4-allyloxy-3,5-dimethoxyphenethylamine (other name: Allylescaline);
 292 4-Bromo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25B-NBOH);
 293 4-chloro-alpha-methylamino-valerophenone (other name: 4-chloropentedrone);
 294 4-chloro-alpha-Pyrrolidinovalerophenone (other name: 4-chloro-alpha-PVP);
 295 4-fluoro-alpha-Pyrrolidinoheptiophenone (other name: 4-fluoro-PV8);
 296 4-hydroxy-N,N-diisopropyltryptamine (other name: 4-OH-DIPT);
 297 4-methyl-alpha-ethylaminopentiophenone;
 298 4-methyl-alpha-Pyrrolidinohexiophenone (other name: MPHP);
 299 5-methoxy-N,N-dimethyltryptamine (other name: 5-MeO-DMT);
 300 5-methoxy-N-ethyl-N-isopropyltryptamine (other name: 5-MeO-EIPT);
 301 6-ethyl-6-nor-lysergic acid diethylamide (other name: ETH-LAD);
 302 6-allyl-6-nor-lysergic acid diethylamide (other name: AL-LAD);
 303 (N-methyl aminopropyl)-2,3-dihydrobenzofuran (other name: MAPDB);
 304 2-(methylamino)-2-phenyl-cyclohexanone (other name: Deschloroketamine);

- 305 2-(ethylamino)-2-phenyl-cyclohexanone (other name: deschloro-N-ethyl-ketamine);
 306 2-methyl-1-(4-(methylthio)phenyl)-2-morpholinopropiophenone (other name: MMMP);
 307 Alpha-ethylaminohexanophenone (other name: N-ethylhexedrone);
 308 N-ethyl-1-(3-methoxyphenyl)cyclohexylamine (other name: 3-methoxy-PCE);
 309 4-fluoro-alpha-pyrrolidinohexiophenone (other name: 4-fluoro-alpha-PHP);
 310 N-ethyl-1,2-diphenylethylamine (other name: Ephedrine);
 311 2,5-dimethoxy-4-chloroamphetamine (other name: DOC);
 312 3,4-methylenedioxy-N-tert-butylcathinone;
 313 Alpha-pyrrolidinoisohexiophenone (other name: alpha-PiHP);
 314 1-[1-(3-hydroxyphenyl)cyclohexyl]piperidine (other name: 3-hydroxy PCP);
 315 4-acetyloxy-N,N-diallyltryptamine (other name: 4-AcO-DALT);
 316 4-hydroxy-N,N-methylisopropyltryptamine (other name: 4-hydroxy-MiPT);
 317 3,4-Methylenedioxy-alpha-pyrrolidinohexanophenone (other name: MDPHP);
 318 5-methoxy-N,N-dibutyltryptamine (other name: 5-methoxy-DBT);
 319 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (other names: Eutylone, bk-EBDB);
 320 1-(1,3-benzodioxol-5-yl)-2-(butylamino)-1-pentanone (other name: N-butylpentylone);
 321 N-benzyl-3,4-dimethoxyamphetamine (other name: N-benzyl-3,4-DMA);
 322 1-(benzo[d][1,3]dioxol-5-yl)-2-(sec-butylamino)pentan-1-one (other name: N-sec-butyl Pentylone);
 323 1-cyclopropionyl lysergic acid diethylamide (other name: 1cP-LSD);
 324 2-(ethylamino)-1-phenylheptan-1-one (other name: N-ethylheptedrone);
 325 (2-ethylaminopropyl)benzofuran (other name: EAPB);
 326 4-ethyl-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25E-NBOH);
 327 2-fluoro-Deschloroketamine (other name: 2-(2-fluorophenyl)-2-(methylamino)-cyclohexanone);
 328 4-hydroxy-N-ethyl-N-propyltryptamine (other name: 4-hydroxy-EPT);
 329 2-(isobutylamino)-1-phenylhexan-1-one (other names: N-Isobutyl Hexedrone,
 330 alpha-isobutylaminohexanphenone);
 331 1-(4-methoxyphenyl)-N-methylpropan-2-amine (other names: para-Methoxymethamphetamine,
 332 PMMA);
 333 N-ethyl-1-(3-hydroxyphenyl)cyclohexylamine (other name: 3-hydroxy-PCE);
 334 N-heptyl-3,4-dimethoxyamphetamine (other names: name: N-heptyl-3,4-DMA);
 335 N-hexyl-3,4-dimethoxyamphetamine (other names: name: N-hexyl-3,4-DMA);
 336 4-fluoro-3-methyl-alpha-pyrrolidinovalerophenone (other name: 4-fluoro-3-methyl-alpha-PVP);
 337 4-fluoro-alpha-methylamino-valerophenone (other name: 4-fluoropentedrone);
 338 N-(1,4-dimethylpentyl)-3,4-dimethoxyamphetamine (other name: N-(1,4-dimethylpentyl)-3,4-DMA);
 339 4,5-methylenedioxy-N,N-diisopropyltryptamine (other name: 4,5-MDO-DiPT);
 340 Alpha-pyrrolidinocyclohexanophenone (other name: alpha-PCYP);
 341 3,4-methylenedioxy-alpha-pyrrolidinoheptiophenone (other name: MDPV8);
 342 4-chloro-alpha-methylaminobutiophenone (other name: 4-chloro Buphedrone).
 343 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 344 or preparation which contains any quantity of the following substances having a depressant effect on the
 345 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such
 346 salts, isomers and salts of isomers is possible within the specific chemical designation:
 347 5-(2-chlorophenyl)-1,3-dihydro-3-methyl-7-nitro-2H-1,4-benzodiazepin-2-one (other name:
 348 Meclonazepam);
 349 7-chloro-5-(2-fluorophenyl)-1,3-dihydro-1,4-benzodiazepin-2-one (other name: Norfludiazepam);
 350 Bromazolam;
 351 Clonazolam;
 352 Deschloroetizolam;
 353 Etizolam;
 354 Flualprazolam;
 355 Flubromazepam;
 356 Flubromazolam;
 357 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;
 358 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
 359 Mecloqualone;
 360 Methaqualone.
 361 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 362 or preparation which contains any quantity of the following substances having a stimulant effect on the
 363 central nervous system, including its salts, isomers and salts of isomers:
 364 2-(3-fluorophenyl)-3-methylmorpholine (other name: 3-fluorophenmetrazine);
 365 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline;
 366 4,5-dihydro-5-phenyl-2-oxazolamine);

367 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,
 368 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;
 369 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazamine);
 370 Ethylamphetamine;
 371 Ethyl phenyl(piperidin-2-yl)acetate (other name: Ethylphenidate);
 372 Fenethylamine;
 373 Methcathinone (some other names: 2-(methylamino)-propionophenone;
 374 alpha-(methylamino)-propionophenone; 2-(methylamino)-1-phenylpropan-1-one;
 375 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;
 376 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);
 377 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
 378 N,N-dimethylamphetamine (other names: N, N-alpha-trimethyl-benzeneethanamine, N,
 379 N-alpha-trimethylphenethylamine);
 380 Methyl 2-(4-fluorophenyl)-2-(2-piperidinyl)acetate (other name: 4-fluoromethylphenidate);
 381 Isopropyl-2-phenyl-2-(2-piperidinyl)acetate (other name: Isopropylphenidate);
 382 4-chloro-N,N-dimethylcathinone;
 383 3,4-methylenedioxy-N-benzylcathinone (other name: BMDP).
 384 6. Any substance that contains one or more cannabimimetic agents or that contains their salts,
 385 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is
 386 possible within the specific chemical designation, and any preparation, mixture, or substance containing,
 387 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.
 388 a. "Cannabimimetic agents" includes any substance that is within any of the following structural
 389 classes:
 390 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or
 391 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;
 392 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of
 393 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not
 394 substituted on the naphthoyl or naphthyl ring to any extent;
 395 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not
 396 further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to
 397 any extent;
 398 1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not
 399 further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to
 400 any extent;
 401 3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring,
 402 whether or not further substituted in the indole ring to any extent, whether or not substituted on the
 403 phenyl ring to any extent;
 404 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring, whether or not further
 405 substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any
 406 extent;
 407 3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further
 408 substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any
 409 extent;
 410 N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring,
 411 whether or not further substituted on the indole ring to any extent, whether or not substituted on the
 412 adamantyl ring to any extent; and
 413 N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring,
 414 whether or not further substituted on the indazole ring to any extent, whether or not substituted on the
 415 adamantyl ring to any extent.
 416 b. The term "cannabimimetic agents" includes:
 417 5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);
 418 5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);
 419 5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);
 420 5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);
 421 1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);
 422 1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);
 423 1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);
 424 1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);
 425 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);
 426 (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tet
 427 rahydrobenzo[c]chromen-1-ol (other name: HU-210);

- 428 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);
- 429 1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);
- 430 1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);
- 431 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);
- 432 1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);
- 433 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);
- 434 1-((N-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);
- 435 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);
- 436 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);
- 437 Pravadoline (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone (other name: WIN 48,098);
- 438 1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);
- 440 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);
- 441 1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);
- 442 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other names: XLR-11, 5-fluoro-UR-144);
- 444 N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);
- 445 N-adamantyl-1-pentylindazole-3-carboxamide (other names: AKB48, APINACA);
- 446 1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);
- 447 (8-quinolinyl)(1-pentylindol-3-yl)carboxylate (other name: PB-22);
- 448 (8-quinolinyl)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);
- 449 (8-quinolinyl)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);
- 450 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);
- 451 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name: AB-FUBINACA);
- 452 1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);
- 454 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: ADB-PINACA);
- 456 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name: AB-CHMINACA);
- 457 AB-CHMINACA);
- 458 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name: 5-fluoro-AB-PINACA);
- 459 5-fluoro-AB-PINACA);
- 460 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other names: ADB-CHMINACA, MAB-CHMINACA);
- 461 names: ADB-CHMINACA, MAB-CHMINACA);
- 462 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name: 5-fluoro-AMB);
- 463 5-fluoro-AMB);
- 464 1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);
- 465 1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);
- 466 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name: MAM-2201);
- 467 N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamide (other name: ADB-FUBINACA);
- 468 (other name: ADB-FUBINACA);
- 469 Methyl 2-[1-(4-fluorophenyl)methyl]-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name: MDMB-FUBINACA);
- 470 name: MDMB-FUBINACA);
- 471 Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names: 5-fluoro-ADB, 5-Fluoro-MDMB-PINACA);
- 472 5-fluoro-ADB, 5-Fluoro-MDMB-PINACA);
- 473 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other names: AMB-FUBINACA, FUB-AMB);
- 474 names: AMB-FUBINACA, FUB-AMB);
- 475 N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other name: 5F-APINACA);
- 476 5F-APINACA);
- 477 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (other name: 5F-AKB48);
- 478 N-(adamantan-1-yl)-1-(5-chloropentyl)indazole-3-carboxamide (other name: 5-chloro-AKB48);
- 479 Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate (other name: SDB-005);
- 480 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide (other name: AB-CHMICA);
- 481 AB-CHMICA);
- 482 1-pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide (other name: SDB-006);
- 483 Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate (other name: FUB-PB-22);
- 484 Methyl N-[1-(cyclohexylmethyl)-1H-indole-3-carbonyl]valinate (other name: MMB-CHMICA);
- 485 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name: 5-fluoro-ADB-PINACA);
- 486 5-fluoro-ADB-PINACA);
- 487 1-(4-cyanobutyl)-N-(1-methyl-1-phenylethyl)-1H-indazole-3-carboxamide (other name: 4-cyano CUMYL-BUTINACA);
- 488 CUMYL-BUTINACA);
- 489 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name: 5-fluoro-ADB-PINACA);

490 ~~5-Fluoro-MDMB-PICA~~) names: *5-fluoro MDMB-PICA, 5F-MDMB-PICA*);
 491 Ethyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other name:
 492 EMB-FUBINACA);
 493 Methyl 2-[1-(4-fluorobutyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name:
 494 4-fluoro-MDMB-BUTINACA);
 495 1-(5-fluoropentyl)-N-(1-methyl-1-phenylethyl)-1H-indole-3-carboxamide (other name: 5-fluoro
 496 CUMYL-PICA);
 497 Methyl 2-[1-(pent-4-enyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name:
 498 MDMB-4en-PINACA);
 499 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indole-3-carbonyl}amino)-3-methylbutanoate (other names:
 500 MMB-FUBICA, AMB-FUBICA);
 501 Methyl 2-[1-(4-penten-1-yl)-1H-indole-3-carboxamido]-3-methylbutanoate (other names: MMB022,
 502 MMB-4en-PICA);
 503 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-methylbutanoate (other name: MMB 2201);
 504 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-phenylpropanoate (other name:
 505 5-fluoro-MPP-PICA);
 506 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butylandazole-3-carboxamide (other name:
 507 ADB-BUTINACA);
 508 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-chloropentyl)indazole-3-carboxamide (other name:
 509 5-chloro-AB-PINACA);
 510 *1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (other names:*
 511 *5F-CUMYL-PINACA, 5-fluoro CUMYL-PINACA, CUMYL-5F-PINACA)*;
 512 *Ethyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names:*
 513 *5F-EDMB-PINACA, 5-fluoro EDMB-PINACA)*;
 514 *Ethyl-2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3-methylbutanoate (other names:*
 515 *5-fluoro-EMB-PINACA, 5F-AEB)*;
 516 *Ethyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-methylbutanoate (other name:*
 517 *5-fluoro-EMB-PICA)*;
 518 *Ethyl-2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name: 5-fluoro*
 519 *EDMB-PICA)*;
 520 *Methyl 2-[1-(4-fluorobutyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name:*
 521 *4-fluoro-MDMB-BUTICA)*;
 522 *Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate (other names:*
 523 *MDMB-CHMICA, MMB-CHMINACA)*;
 524 *N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-enyl)indazole-3-carboxamide (other name:*
 525 *ADB-4en-PINACA)*.