

VIRGINIA ACTS OF ASSEMBLY — CHAPTER

An Act to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Schedule I controlled substances.

[H 1194]

Approved

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

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

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-alphacetylmethadol, 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;

Levophenacymorphan;

Morpheridine;

N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (other name: Cyclopropyl fentanyl);

N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (other name: Tetrahydrofuran fentanyl);

N-[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]-N-phenylpropanamide (other name: alpha-methylthiofentanyl);

N-[1-(1-methyl-2-phenylethyl)-4-piperidyl]-N-phenylacetamide (other name: acetyl-alpha-methylfentanyl);

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- 57 *N*-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyll}-*N*-phenylpropanamide (other name:
 58 *beta*-hydroxythiofentanyl);
 59 *N*-[1-(2-hydroxy-2-phenyl)ethyl-4-piperidyl]-*N*-phenylpropanamide (other name:
 60 *beta*-hydroxyfentanyl);
 61 *N*-[1-(α -methyl-*beta*-phenyl)ethyl-4-piperidyl]propionanilide (other names:
 62 1-(1-methyl-2-phenylethyl)-4-(*N*-propanilido) piperidine, α -methylfentanyl);
 63 *N*-(2-fluorophenyl)-*N*-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other names: 2-fluorofentanyl,
 64 *ortho*-fluorofentanyl);
 65 *N*-(3-fluorophenyl)-*N*-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name: 3-fluorofentanyl);
 66 *N*-[3-methyl-1-(2-hydroxy-2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide (other name:
 67 *beta*-hydroxy-3-methylfentanyl);
 68 *N*-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide (other name: 3-methylfentanyl);
 69 *N*-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyll]-*N*-phenylpropanamide (other name:
 70 3-methylthiofentanyl);
 71 *N*-(4-fluorophenyl)-2-methyl-*N*-[1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name:
 72 *para*-fluoroisobutyryl fentanyl);
 73 *N*-(4-fluorophenyl)-*N*-[1-(2-phenylethyl)-4-piperidinyll]-butanamide (other name:
 74 *para*-fluorobutyrylfentanyl);
 75 *N*-(4-fluorophenyl)-*N*-1-(2-phenylethyl)-4-piperidinyll]-propanamide (other name: *para*-fluorofentanyl);
 76 Noracymethadol;
 77 Norlevorphanol;
 78 Normethadone;
 79 Norpipanone;
 80 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyll]-2-furancarboxamide (other name: Furanyl fentanyl);
 81 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyll]-2-propenamide (other name: Acryl fentanyl);
 82 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyll]-butanamide (other name: butyryl fentanyl);
 83 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyll]-pentanamide (other name: Pentanoyl fentanyl);
 84 *N*-phenyl-*N*-[1-(2-thienyl)ethyl-4-piperidinyll]-propanamide (other name: thiofentanyl);
 85 Phenadoxone;
 86 Phenampromide;
 87 Phenomorphan;
 88 Phenoperidine;
 89 Piritramide;
 90 Proheptazine;
 91 Properidine;
 92 Propiram;
 93 Racemoramide;
 94 Tilidine;
 95 Trimeperidine.
- 96 2. Any of the following opium derivatives, their salts, isomers and salts of isomers, unless
 97 specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible
 98 within the specific chemical designation:
 99 Acetorphine;
 100 Acetyldihydrocodeine;
 101 Benzylmorphine;
 102 Codeine methylbromide;
 103 Codeine-*N*-Oxide;
 104 Cyprenorphine;
 105 Desomorphine;
 106 Dihydromorphine;
 107 Drotebanol;
 108 Etorphine;
 109 Heroin;
 110 Hydromorphanol;
 111 Methyldesorphine;
 112 Methyldihydromorphine;
 113 Morphine methylbromide;
 114 Morphine methylsulfonate;
 115 Morphine-*N*-Oxide;
 116 Myrophine;
 117 Nicocodeine;

118 Nicomorphine;
 119 Normorphine;
 120 Pholcodine;
 121 Thebacon.
 122 3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture,
 123 or preparation, which contains any quantity of the following hallucinogenic substances, or which
 124 contains any of its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers,
 125 and salts of isomers is possible within the specific chemical designation (for purposes of this subdivision
 126 only, the term "isomer" includes the optical, position, and geometric isomers):
 127 Alpha-ethyltryptamine (some trade or other names: Monase; a-ethyl-1H-indole-3-ethanamine;
 128 3-2-aminobutyl] indole; a-ET; AET);
 129 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:
 130 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B; Nexus);
 131 3,4-methylenedioxy amphetamine;
 132 5-methoxy-3,4-methylenedioxy amphetamine;
 133 3,4,5-trimethoxy amphetamine;
 134 Alpha-methyltryptamine (other name: AMT);
 135 Bufotenine;
 136 Diethyltryptamine;
 137 Dimethyltryptamine;
 138 4-methyl-2,5-dimethoxyamphetamine;
 139 2,5-dimethoxy-4-ethylamphetamine (DOET);
 140 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
 141 Ibogaine;
 142 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
 143 Lysergic acid diethylamide;
 144 Mescaline;
 145 Parahexyl (some trade or other names:
 146 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl);
 147 Peyote;
 148 N-ethyl-3-piperidyl benzilate;
 149 N-methyl-3-piperidyl benzilate;
 150 Psilocybin;
 151 Psilocyn;
 152 Salvinorin A;
 153 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated
 154 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;
 155 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish);
 156 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;
 157 2,5-DMA);
 158 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts
 159 and salts of isomers;
 160 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4
 161 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);
 162 N-hydroxy-3,4-methylenedioxyamphetamine (some other names:
 163 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);
 164 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:
 165 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);
 166 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;
 167 paramethoxyamphetamine; PMA);
 168 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,
 169 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);
 170 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy,
 171 PHP);
 172 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl) -cyclohexyl]-piperidine,
 173 2-thienyl analog of phencyclidine, TPCP, TCP);
 174 1-1-(2-thienyl)cyclohexyl]pyrrolidine (other name: TCPy);
 175 3,4-methylenedioxypyrovalerone (other name: MDPV);
 176 4-methylmethcathinone (other names: mephedrone, 4-MMC);
 177 3,4-methylenedioxymethcathinone (other name: methylone);
 178 Naphthylpyrovalerone (other name: naphyrone);

- 179 4-fluoromethcathinone (other name: flephedrone, 4-FMC);
 180 4-methoxymethcathinone (other names: methedrone; bk-PMMA);
 181 Ethcathinone (other name: N-ethylcathinone);
 182 3,4-methylenedioxyethcathinone (other name: ethylone);
 183 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone);
 184 N,N-dimethylcathinone (other name: metamfepramone);
 185 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);
 186 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);
 187 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);
 188 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);
 189 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);
 190 3-fluoromethcathinone (other name: 3-FMC);
 191 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);
 192 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);
 193 4-Methylethcathinone (other name: 4-MEC);
 194 4-Ethylmethcathinone (other name: 4-EMC);
 195 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);
 196 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP);
 197 Alpha-methylamino-butyrophenone (other name: Buphedrone);
 198 Alpha-methylamino-valerophenone (other name: Pentedrone);
 199 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);
 200 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);
 201 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,
 202 25I-NBOMe, 2C-I-NBOMe);
 203 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);
 204 4-Fluoromethamphetamine (other name: 4-FMA);
 205 4-Fluoroamphetamine (other name: 4-FA);
 206 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);
 207 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);
 208 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);
 209 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);
 210 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);
 211 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
 212 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
 213 (2-aminopropyl)benzofuran (other name: APB);
 214 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
 215 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 216 2C-C-NBOMe, 25C-NBOMe, 25C);
 217 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 218 2C-B-NBOMe, 25B-NBOMe, 25B);
 219 Acetoxydimethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
 220 Benocyclidine (other names: BCP, BTCP);
 221 Alpha-pyrrolidinobutiophenone (other name: alpha-PBP);
 222 3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
 223 4-bromomethcathinone (other name: 4-BMC);
 224 4-chloromethcathinone (other name: 4-CMC);
 225 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
 226 Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
 227 Alpha-Pyrrolidinoheptiophenone (other name: PV8);
 228 5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
 229 Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
 230 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
 231 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
 232 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
 233 4-Chloroethcathinone (other name: 4-CEC);
 234 3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
 235 1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
 236 (2-Methylaminopropyl)benzofuran (other name: MAPB);
 237 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)-1-pentanone (other names: N,N-Dimethylpentylone,
 238 Dipentylone);
 239 1-(4-methoxyphenyl)-2-(pyrrolidin-1-yl)octan-1-one (other name: 4-methoxy-PV9)

- 240 3,4-tetramethylene-alpha-pyrrolidinovalerophenone (other name: TH-PVP);
 241 4-allyloxy-3,5-dimethoxyphenethylamine (other name: Allylescaline);
 242 4-Bromo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25B-NBOH);
 243 4-chloro-alpha-methylamino-valerophenone (other name: 4-chloropentedrone);
 244 4-chloro-alpha-Pyrrolidinovalerophenone (other name: 4-chloro-alpha-PVP);
 245 4-fluoro-alpha-Pyrrolidinoheptiophenone (other name: 4-fluoro-PV8);
 246 4-hydroxy-N,N-diisopropyltryptamine (other name: 4-OH-DIPT);
 247 4-methyl-alpha-ethylaminopentiophenone;
 248 4-methyl-alpha-Pyrrolidinohexiophenone (other name: MPHP);
 249 5-methoxy-N,N-dimethyltryptamine (other name: 5-MeO-DMT);
 250 5-methoxy-N-ethyl-N-isopropyltryptamine (other name: 5-MeO-EIPT);
 251 6-ethyl-6-nor-lysergic acid diethylamide (other name: ETH-LAD);
 252 6-allyl-6-nor-lysergic acid diethylamide (other name: AL-LAD);
 253 (N-methyl aminopropyl)-2,3-dihydrobenzofuran (other name: MAPDB).
 254 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 255 or preparation which contains any quantity of the following substances having a depressant effect on the
 256 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such
 257 salts, isomers and salts of isomers is possible within the specific chemical designation:
 258 Clonazepam;
 259 Etizolam;
 260 Flubromazepam;
 261 Flubromazolam;
 262 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;
 263 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
 264 Mecloqualone;
 265 Methaqualone.
 266 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 267 or preparation which contains any quantity of the following substances having a stimulant effect on the
 268 central nervous system, including its salts, isomers and salts of isomers:
 269 2-(3-fluorophenyl)-3-methylmorpholine (other name: 3-fluorophenmetrazine);
 270 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline;
 271 4,5-dihydro-5-phenyl-2-oxazolamine);
 272 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,
 273 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;
 274 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
 275 Ethylamphetamine;
 276 Ethyl phenyl(piperidin-2-yl)acetate (other name: Ethylphenidate);
 277 Fenethylamine;
 278 Methcathinone (some other names: 2-(methylamino)-propionophenone;
 279 alpha-(methylamino)-propionophenone; 2-(methylamino)-1-phenylpropan-1-one;
 280 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;
 281 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);
 282 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
 283 N,N-dimethylamphetamine (other names: N, N-alpha-trimethyl-benzeneethanamine, N,
 284 N-alpha-trimethylphenethylamine).
 285 6. Any substance that contains one or more cannabimimetic agents or that contains their salts,
 286 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is
 287 possible within the specific chemical designation, and any preparation, mixture, or substance containing,
 288 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.
 289 a. "Cannabimimetic agents" includes any substance that is within any of the following structural
 290 classes:
 291 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or
 292 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;
 293 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of
 294 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not
 295 substituted on the naphthoyl or naphthyl ring to any extent;
 296 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not
 297 further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to
 298 any extent;
 299 1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not
 300 further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to

301 any extent;
 302 3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring,
 303 whether or not further substituted in the indole ring to any extent, whether or not substituted on the
 304 phenyl ring to any extent;
 305 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring, whether or not further
 306 substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any
 307 extent;
 308 3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further
 309 substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any
 310 extent;
 311 N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring,
 312 whether or not further substituted on the indole ring to any extent, whether or not substituted on the
 313 adamantyl ring to any extent; and
 314 N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring,
 315 whether or not further substituted on the indazole ring to any extent, whether or not substituted on the
 316 adamantyl ring to any extent.
 317 b. The term "cannabimimetic agents" includes:
 318 5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);
 319 5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);
 320 5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);
 321 5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);
 322 1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);
 323 1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);
 324 1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);
 325 1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);
 326 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);
 327 (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-terahydrobenzo[c]chr-
 328 omen-1-ol (other name: HU-210);
 329 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);
 330 1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);
 331 1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);
 332 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);
 333 1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);
 334 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);
 335 1-((N-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);
 336 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);
 337 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);
 338 Pravadoline (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone (other
 339 name: WIN 48,098);
 340 1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);
 341 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);
 342 1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);
 343 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other names: XLR-11,
 344 5-fluoro-UR-144);
 345 N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);
 346 N-adamantyl-1-pentylindazole-3-carboxamide (other names: AKB48, APINACA);
 347 1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);
 348 (8-quinoliny)(1-pentylindol-3-yl)carboxylate (other name: PB-22);
 349 (8-quinoliny)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);
 350 (8-quinoliny)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);
 351 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);
 352 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name:
 353 AB-FUBINACA);
 354 1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);
 355 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name:
 356 ADB-PINACA);
 357 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name:
 358 AB-CHMINACA);
 359 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:
 360 5-fluoro-AB-PINACA);
 361 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other names:

362 ADB-CHMINACA, MAB-CHMINACA);
 363 Methyl-2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name:
 364 5-fluoro-AMB);
 365 1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);
 366 1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);
 367 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name MAM-2201);
 368 N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamide
 369 (other name: ADB-FUBINACA);
 370 Methyl 2-[1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name:
 371 MDMB-FUBINACA);
 372 Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names:
 373 5-fluoro-ADB, 5-Fluoro-MDMB-PINACA);
 374 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other
 375 names: AMB-FUBINACA, FUB-AMB);
 376 N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other name: FUB-AKB48)
 377 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (other name: 5F-AKB48);
 378 Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate (other name: SDB-005);
 379 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indole -3-carboxamide (other name:
 380 AB-CHMICA);
 381 1-pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide (other name: SDB-006);
 382 Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate (other name: FUB-PB-22);
 383 Methyl N-[1-(cyclohexylmethyl)-1H-indole-3-carbonyl]valinate (other name: MMB-CHMICA);
 384 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:
 385 5-fluoro-ADB-PINACA).
 386 **2. That the provisions of this act may result in a net increase in periods of imprisonment or**
 387 **commitment. Pursuant to § 30-19.1:4, the estimated amount of the necessary appropriation is \$0**
 388 **for periods of imprisonment in state adult correctional facilities and cannot be determined for**
 389 **periods of commitment to the custody of the Department of Juvenile Justice.**