

VIRGINIA ACTS OF ASSEMBLY — CHAPTER

An Act to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Drug Control Act; Schedule I drugs; addition of substances.

[S 1546]

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);

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-acetylmethadol, 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;

Etoxadine;

Furethidine;

Hydroxypethidine;

Ketobemidone;

Levomoramide;

Levophenacetylmorphan;

Morpheridine;

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);

N-[1-(2-hydroxy-2-phenyl)ethyl-4-piperidyl]-N-phenylpropanamide (other name: beta-hydroxyfentanyl);

N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide (other names: 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine, alpha-methylfentanyl);

N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: 3-fluorofentanyl);

57 *N*-[3-methyl-1-(2-hydroxy-2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide (other name:
 58 *beta*-hydroxy-3-methylfentanyl);
 59 *N*-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide (other name: 3-methylfentanyl);
 60 *N*-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-*N*-phenylpropanamide (other name:
 61 3-methylthiofentanyl);
 62 *N*-(4-fluorophenyl)-*N*-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name:
 63 *para*-fluorobutyrylfentanyl);
 64 *N*-(4-fluorophenyl)-*N*-1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: *para*-fluorofentanyl);
 65 Noracymethadol;
 66 Norlevorphanol;
 67 Normethadone;
 68 Norpipanone;
 69 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyl]-2-furancarboxamide (other name: *Furanyl fentanyl*);
 70 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: *butyryl fentanyl*);
 71 *N*-phenyl-*N*-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide (other name: *Pentanoyl fentanyl*);
 72 *N*-phenyl-*N*-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide (other name: *thiofentanyl*);
 73 Phenadoxone;
 74 Phenampromide;
 75 Phenomorphan;
 76 Phenoperidine;
 77 Piritramide;
 78 Proheptazine;
 79 Properidine;
 80 Propiram;
 81 Racemoramide;
 82 Tilidine;
 83 Trimeperidine.

84 2. Any of the following opium derivatives, their salts, isomers and salts of isomers, unless
 85 specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible
 86 within the specific chemical designation:

87 Acetorphine;
 88 Acetyldihydrocodeine;
 89 Benzylmorphine;
 90 Codeine methylbromide;
 91 Codeine-*N*-Oxide;
 92 Cyprenorphine;
 93 Desomorphine;
 94 Dihydromorphine;
 95 Drotebanol;
 96 Etorphine;
 97 Heroin;
 98 Hydromorphanol;
 99 Methyldesorphine;
 100 Methyldihydromorphine;
 101 Morphine methylbromide;
 102 Morphine methylsulfonate;
 103 Morphine-*N*-Oxide;
 104 Myrophine;
 105 Nicocodeine;
 106 Nicomorphine;
 107 Normorphine;
 108 Pholcodine;
 109 Thebacon.

110 3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture,
 111 or preparation, which contains any quantity of the following hallucinogenic substances, or which
 112 contains any of its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers,
 113 and salts of isomers is possible within the specific chemical designation (for purposes of this subdivision
 114 only, the term "isomer" includes the optical, position, and geometric isomers):

115 Alpha-ethyltryptamine (some trade or other names: Monase; α -ethyl-1H-indole-3-ethanamine;
 116 3-2-aminobutyl] indole; α -ET; AET);
 117 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:

- 118 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);
 119 3,4-methylenedioxy amphetamine;
 120 5-methoxy-3,4-methylenedioxy amphetamine;
 121 3,4,5-trimethoxy amphetamine;
 122 Alpha-methyltryptamine (other name: AMT);
 123 Bufotenine;
 124 Diethyltryptamine;
 125 Dimethyltryptamine;
 126 4-methyl-2,5-dimethoxyamphetamine;
 127 2,5-dimethoxy-4-ethylamphetamine (DOET);
 128 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
 129 Ibogaine;
 130 5-methoxy-N, N-diisopropyltryptamine (other name: 5-MeO-DIPT);
 131 Lysergic acid diethylamide;
 132 Mescaline;
 133 Parahexyl (some trade or other names:
 134 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo -b, d] pyran; Synhexyl);
 135 Peyote;
 136 N-ethyl-3-piperidyl benzilate;
 137 N-methyl-3-piperidyl benzilate;
 138 Psilocybin;
 139 Psilocyn;
 140 Salvinorin A;
 141 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated
 142 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;
 143 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish);
 144 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;
 145 2,5-DMA);
 146 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts
 147 and salts of isomers;
 148 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4
 149 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);
 150 N-hydroxy-3,4-methylenedioxyamphetamine (some other names:
 151 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);
 152 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:
 153 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);
 154 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;
 155 paramethoxyamphetamine; PMA);
 156 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,
 157 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);
 158 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy,
 159 PHP);
 160 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl) -cyclohexyl]-piperidine,
 161 2-thienyl analog of phencyclidine, TPCP, TCP);
 162 1-1-(2-thienyl)cyclohexyl]pyrrolidine (other name: TCPy);
 163 3,4-methylenedioxypyrovalerone (other name: MDPV);
 164 4-methylmethcathinone (other names: mephedrone, 4-MMC);
 165 3,4-methylenedioxymethcathinone (other name: methylone);
 166 Naphthylpyrovalerone (other name: naphyrone);
 167 4-fluoromethcathinone (other name: flephedrone, 4-FMC);
 168 4-methoxymethcathinone (other names: methedrone; bk-PMMA);
 169 Ethcathinone (other name: N-ethylcathinone);
 170 3,4-methylenedioxyethcathinone (other name: ethylone);
 171 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine benzodioxolylbutanamine (other name: butylone);
 172 N, N-dimethylcathinone (other name: metamfepramone);
 173 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);
 174 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);
 175 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);
 176 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);
 177 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);
 178 3-fluoromethcathinone (other name: 3-FMC);

- 179 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);
 180 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);
 181 4-Methylethcathinone (other name: 4-MEC);
 182 4-Ethylmethcathinone (other name: 4-EMC);
 183 N, N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);
 184 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP);
 185 Alpha-methylamino-butyrophenone (other name: Buphedrone);
 186 Alpha-methylamino-valerophenone (other name: Pentedrone);
 187 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);
 188 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);
 189 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,
 190 25I-NBOMe, 2C-I-NBOMe);
 191 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);
 192 4-Fluoromethamphetamine (other name: 4-FMA);
 193 4-Fluoroamphetamine (other name: 4-FA);
 194 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);
 195 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);
 196 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);
 197 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);
 198 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);
 199 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
 200 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
 201 (2-aminopropyl)benzofuran (other name: APB);
 202 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
 203 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 204 2C-C-NBOMe, 25C-NBOMe, 25C);
 205 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 206 2C-B-NBOMe, 25B-NBOMe, 25B);
 207 Acetoxymethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
 208 Benocyclidine (other names: BCP, BTCP);
 209 Alpha-pyrrolidinobutiophenone (other name: alpha-PBP);
 210 3,4-methylenedioxy-N, N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
 211 4-bromomethcathinone (other name: 4-BMC);
 212 4-chloromethcathinone (other name: 4-CMC);
 213 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
 214 Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
 215 Alpha-Pyrrolidinoheptiophenone (other name: PV8);
 216 5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
 217 Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
 218 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
 219 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
 220 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
 221 4-Chloroethcathinone (other name: 4-CEC);
 222 3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
 223 1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
 224 (2-Methylaminopropyl)benzofuran (other name: MAPB).
 225 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 226 or preparation which contains any quantity of the following substances having a depressant effect on the
 227 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such
 228 salts, isomers and salts of isomers is possible within the specific chemical designation:
 229 Clonazepam;
 230 Etizolam;
 231 Flubromazepam;
 232 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;
 233 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
 234 Mecloqualone;
 235 Methaqualone.
 236 ~~Etizolam.~~
 237 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 238 or preparation which contains any quantity of the following substances having a stimulant effect on the
 239 central nervous system, including its salts, isomers and salts of isomers:

- 240 2-(3-fluorophenyl)-3-methylmorpholine (other name: 3-fluorophenmetrazine);
 241 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline;
 242 4,5-dihydro-5-phenyl-2-oxazolamine);
 243 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
 244 Fenethylline;
 245 Ethylamphetamine;
 246 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,
 247 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;
 248 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
 249 Ethylamphetamine;
 250 Ethyl phenyl(piperidin-2-yl)acetate (other name: Ethylphenidate);
 251 Fenethylline;
 252 Methcathinone (some other names: 2-(methylamino)-propionophenone;
 253 alpha-(methylamino)-propionophenone; 2-(methylamino)-1-phenylpropan-1-one;
 254 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;
 255 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);
 256 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
 257 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
 258 N, N-dimethylamphetamine (other names: N, N-alpha-trimethyl-benzeneethanamine, N,
 259 N-alpha-trimethylphenethylamine).
 260 6. Any material, compound, mixture or preparation containing any quantity of the following
 261 substances:
 262 N-3-methyl-1-(2-phenethyl)-4-piperidyl]-N-phenylpropanamide (other name: 3-methylfentanyl), its
 263 optical and geometric isomers, salts, and salts of isomers;
 264 1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP), its optical isomers, salts and salts of
 265 isomers;
 266 1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP), its optical isomers, salts and
 267 salts of isomers;
 268 N-1-(alpha-methyl-beta-phenyl)-ethyl-4-piperidyl]-propionanilide (other names:
 269 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine), alpha-methylfentanyl);
 270 N-1-(1-methyl-2-phenethyl)-4-piperidyl]-N-phenylacetamide (other name:
 271 acetyl-alpha-methylfentanyl), its optical isomers, salts and salts of isomers;
 272 N-1-(1-methyl-2-2-thienyl)ethyl-4-piperidyl]-N-phenylpropanamide (other name:
 273 alpha-methylthiofentanyl), its optical isomers, salts and salts of isomers;
 274 N-1-benzyl-4-piperidyl]-N-phenylpropanamide (other name: benzylfentanyl), its optical isomers, salts
 275 and salts of isomers;
 276 N-1-(2-hydroxy-2-phenyl)-ethyl-4-piperidyl]-N-phenylpropanamide (other name:
 277 beta-hydroxyfentanyl), its optical isomers, salts and salts of isomers;
 278 N-3-methyl-1-(2-hydroxy-2-phenethyl)-4-piperidyl]-N-phenylpropanamide (other name:
 279 beta-hydroxy-3-methylfentanyl), its optical and geometric isomers, salts and salts of isomers;
 280 N-(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (other name:
 281 3-methylthiofentanyl), its optical and geometric isomers, salts and salts of isomers;
 282 N-1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (other name: thienylfentanyl), its optical
 283 isomers, salts and salts of isomers;
 284 N-phenyl-N-1-(2-thienyl)ethyl-4-piperidinyl]-propanamide (other name: thiofentanyl), its optical
 285 isomers, salts and salts of isomers;
 286 N-(4-fluorophenyl)-N-1-(2-phenethyl)-4-piperidinyl]-propanamide (other name: para-fluorofentanyl),
 287 its optical isomers, salts and salts of isomers;
 288 Acetyl fentanyl (other name: desmethyl fentanyl).
 289 7. Any substance that contains one or more cannabimimetic agents or that contains their salts,
 290 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is
 291 possible within the specific chemical designation, and any preparation, mixture, or substance containing,
 292 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.
 293 a. "Cannabimimetic agents" includes any substance that is within any of the following structural
 294 classes:
 295 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or
 296 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;
 297 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of
 298 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not
 299 substituted on the naphthoyl or naphthyl ring to any extent;
 300 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not

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

362 AB-CHMINACA);
 363 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:
 364 5-fluoro-AB-PINACA);
 365 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other names:
 366 ADB-CHMINACA, MAB-CHMINACA);
 367 Methyl-2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name:
 368 5-fluoro-AMB);
 369 1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);
 370 1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);
 371 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name MAM-2201);
 372 *N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamide*
 373 *(other name: ADB-FUBINACA);*
 374 *Methyl 2-[1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other*
 375 *name: MDMB-FUBINACA);*
 376 *Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names:*
 377 *5-fluoro-ADB, 5-Fluoro-MDMB-PINACA;*
 378 *Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other*
 379 *names: AMB-FUBINACA, FUB-AMB);*
 380 *N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other name: FUB-AKB48)*
 381 *N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (other name: 5F-AKB48);*
 382 *Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate (other name: SDB-005);*
 383 *N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide (other name:*
 384 *AB-CHMICA).*