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HOUSE BILL NO. 1610

House Amendments in [] - January 19, 2017

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

Patron Prior to Engrossment—Delegate Garrett

Referred to Committee on Health, Welfare and Institutions

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-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;

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);

ENGROSSED

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

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

- 90 Acetorphine;
91 Acetyldihydrocodeine;
92 Benzylmorphine;
93 Codeine methylbromide;
94 Codeine-N-Oxide;
95 Cyprenorphine;
96 Desomorphine;
97 Dihydromorphine;
98 Drotebanol;
99 Etorphine;
100 Heroin;
101 Hydromorphanol;
102 Methyl-desorphine;
103 Methyl-dihydromorphine;
104 Morphine methylbromide;
105 Morphine methylsulfonate;
106 Morphine-N-Oxide;
107 Myrophine;
108 Nicocodeine;
109 Nicomorphine;
110 Normorphine;
111 Pholcodine;
112 Thebacon.

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

- 118 Alpha-ethyltryptamine (some trade or other names: Monase; α -ethyl-1H-indole-3-ethanamine;
119 3-2-aminobutyl] indole; α -ET; AET);
120 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:

- 121 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);
 122 3,4-methylenedioxy amphetamine;
 123 5-methoxy-3,4-methylenedioxy amphetamine;
 124 3,4,5-trimethoxy amphetamine;
 125 Alpha-methyltryptamine (other name: AMT);
 126 Bufotenine;
 127 Diethyltryptamine;
 128 Dimethyltryptamine;
 129 4-methyl-2,5-dimethoxyamphetamine;
 130 2,5-dimethoxy-4-ethylamphetamine (DOET);
 131 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
 132 Ibogaine;
 133 5-methoxy-N, N-diisopropyltryptamine (other name: 5-MeO-DIPT);
 134 Lysergic acid diethylamide;
 135 Mescaline;
 136 Parahexyl (some trade or other names:
 137 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo -b, d] pyran; Synhexyl);
 138 Peyote;
 139 N-ethyl-3-piperidyl benzilate;
 140 N-methyl-3-piperidyl benzilate;
 141 Psilocybin;
 142 Psilocyn;
 143 Salvinorin A;
 144 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated
 145 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;
 146 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish);
 147 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;
 148 2,5-DMA);
 149 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts
 150 and salts of isomers;
 151 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4
 152 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);
 153 N-hydroxy-3,4-methylenedioxyamphetamine (some other names:
 154 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);
 155 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:
 156 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);
 157 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;
 158 paramethoxyamphetamine; PMA);
 159 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,
 160 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);
 161 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy,
 162 PHP);
 163 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl) -cyclohexyl]-piperidine,
 164 2-thienyl analog of phencyclidine, TPCP, TCP);
 165 1-1-(2-thienyl)cyclohexyl]pyrrolidine (other name: TCPy);
 166 3,4-methylenedioxypyrovalerone (other name: MDPV);
 167 4-methylmethcathinone (other names: mephedrone, 4-MMC);
 168 3,4-methylenedioxymethcathinone (other name: methylone);
 169 Naphthylpyrovalerone (other name: naphyrone);
 170 4-fluoromethcathinone (other name: flephedrone, 4-FMC);
 171 4-methoxymethcathinone (other names: methedrone; bk-PMMA);
 172 Ethcathinone (other name: N-ethylcathinone);
 173 3,4-methylenedioxyethcathinone (other name: ethylone);
 174 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine benzodioxolylbutanamine (other name: butylone);
 175 N, N-dimethylcathinone (other name: metamfepramone);
 176 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);
 177 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);
 178 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);
 179 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);
 180 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);
 181 3-fluoromethcathinone (other name: 3-FMC);

- 182 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);
 183 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);
 184 4-Methylethcathinone (other name: 4-MEC);
 185 4-Ethylmethcathinone (other name: 4-EMC);
 186 N, N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);
 187 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP);
 188 Alpha-methylamino-butyrophenone (other name: Buphedrone);
 189 Alpha-methylamino-valerophenone (other name: Pentedrone);
 190 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);
 191 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);
 192 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,
 193 25I-NBOMe [, 2C-I-NBOMe]);
 194 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);
 195 4-Fluoromethamphetamine (other name: 4-FMA);
 196 4-Fluoroamphetamine (other name: 4-FA);
 197 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);
 198 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);
 199 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);
 200 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);
 201 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);
 202 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
 203 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
 204 (2-aminopropyl)benzofuran (other name: APB);
 205 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
 206 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 207 2C-C-NBOMe, 25C-NBOMe [, 25C]);
 208 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
 209 2C-B-NBOMe, 25B-NBOMe [, 25B]);
 210 Acetoxydimethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
 211 Benocyclidine (other names: BCP, BTCP);
 212 Alpha-pyrrolidinobutiophenone (other name: alpha-PBP);
 213 3,4-methylenedioxy-N, N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
 214 4-bromomethcathinone (other name: 4-BMC);
 215 4-chloromethcathinone (other name: 4-CMC);
 216 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
 217 Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
 218 Alpha-Pyrrolidinoheptiophenone (other name: PV8);
 219 5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
 220 Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
 221 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
 222 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
 223 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
 224 4-Chloroethcathinone (other name: 4-CEC);
 225 3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
 226 1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
 227 (2-Methylaminopropyl)benzofuran (other name: MAPB).
 228 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 229 or preparation which contains any quantity of the following substances having a depressant effect on the
 230 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such
 231 salts, isomers and salts of isomers is possible within the specific chemical designation:
 232 Clonazepam;
 233 Etizolam;
 234 Flubromazepam;
 235 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;
 236 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
 237 Mecloqualone;
 238 Methaqualone.
 239 ~~Etizolam.~~
 240 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture
 241 or preparation which contains any quantity of the following substances having a stimulant effect on the
 242 central nervous system, including its salts, isomers and salts of isomers:
 243 2-(3-fluorophenyl)-3-methylmorpholine (other name: 3-fluorophenmetrazine);

- 244 Aminorex (some trade or other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline;
 245 4,5-dihydro-5-phenyl-2-oxazamine);
 246 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
 247 Fenethylamine;
 248 Ethylamphetamine;
 249 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,
 250 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;
 251 *Cis-4-methylaminorex* (other name: *cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazamine*);
 252 *Ethylamphetamine*;
 253 *Ethyl phenyl(piperidin-2-yl)acetate* (other name: *Ethylphenidate*);
 254 *Fenethylamine*;
 255 Methcathinone (some other names: 2-(methylamino)-propiofenone;
 256 alpha-(methylamino)-propiofenone; 2-(methylamino)-1-phenylpropan-1-one;
 257 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;
 258 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);
 259 *Cis-4-methylaminorex* (other name: *cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazamine*);
 260 *N-Benzylpiperazine* (some other names: BZP, 1-benzylpiperazine);
 261 N, N-dimethylamphetamine (other names: N, N-alpha-trimethyl-benzeneethanamine, N,
 262 N-alpha-trimethylphenethylamine).
 263 6. Any material, compound, mixture or preparation containing any quantity of the following
 264 substances:
 265 N-3-methyl-1-(2-phenethyl)-4-piperidyl-N-phenylpropanamide (other name: 3-methylfentanyl), its
 266 optical and geometric isomers, salts, and salts of isomers;
 267 1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP), its optical isomers, salts and salts of
 268 isomers;
 269 1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP), its optical isomers, salts and
 270 salts of isomers;
 271 N-1-(alpha-methyl-beta-phenyl)-ethyl-4-piperidyl-propionanilide (other names:
 272 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine), alpha-methylfentanyl);
 273 N-1-(1-methyl-2-phenethyl)-4-piperidyl-N-phenylacetamide (other name:
 274 acetyl-alpha-methylfentanyl), its optical isomers, salts and salts of isomers;
 275 N-1-(1-methyl-2-2-thienyl)ethyl-4-piperidyl-N-phenylpropanamide (other name:
 276 alpha-methylthiofentanyl), its optical isomers, salts and salts of isomers;
 277 N-1-benzyl-4-piperidyl-N-phenylpropanamide (other name: benzylfentanyl), its optical isomers, salts
 278 and salts of isomers;
 279 N-1-(2-hydroxy-2-phenyl)-ethyl-4-piperidyl-N-phenylpropanamide (other name:
 280 beta-hydroxyfentanyl), its optical isomers, salts and salts of isomers;
 281 N-3-methyl-1-(2-hydroxy-2-phenethyl)-4-piperidyl-N-phenylpropanamide (other name:
 282 beta-hydroxy-3-methylfentanyl), its optical and geometric isomers, salts and salts of isomers;
 283 N-(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl)-N-phenylpropanamide (other name:
 284 3-methylthiofentanyl), its optical and geometric isomers, salts and salts of isomers;
 285 N-1-(2-thienyl)methyl-4-piperidyl-N-phenylpropanamide (other name: thienylfentanyl), its optical
 286 isomers, salts and salts of isomers;
 287 N-phenyl-N-1-(2-thienyl)ethyl-4-piperidinyl-propanamide (other name: thiofentanyl), its optical
 288 isomers, salts and salts of isomers;
 289 N-(4-fluorophenyl)-N-1-(2-phenethyl)-4-piperidinyl-propanamide (other name: para-fluorofentanyl),
 290 its optical isomers, salts and salts of isomers;
 291 Acetyl fentanyl (other name: desmethyl fentanyl).
 292 7. Any substance that contains one or more cannabimimetic agents or that contains their salts,
 293 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is
 294 possible within the specific chemical designation, and any preparation, mixture, or substance containing,
 295 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.
 296 a. "Cannabimimetic agents" includes any substance that is within any of the following structural
 297 classes:
 298 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or
 299 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;
 300 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of
 301 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not
 302 substituted on the naphthoyl or naphthyl ring to any extent;
 303 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not
 304 further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to

any extent;

1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to any extent;

3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted in the indole ring to any extent, whether or not substituted on the phenyl ring to any extent;

3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any extent;

3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any extent;

N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any extent; and

N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the adamantyl ring to any extent.

b. The term "cannabimimetic agents" includes:

5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);

5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);

5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);

5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);

1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);

1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);

1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);

1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);

1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);

(6aR, 10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7,10,10a-tetrahydrobenzo[c]chromen-1-ol (other name: HU-210);

1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);

1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);

1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);

1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);

1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);

1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);

1-((N-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);

1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);

1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);

Pravadoline (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone (other name: WIN 48,098);

1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);

1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);

1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);

1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: XLR-11 [5-fluoro-UR-144]);

N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);

N-adamantyl-1-pentylindazole-3-carboxamide (other name: AKB48 [, APINACA]);

1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);

(8-quinoliny)(1-pentylindol-3-yl)carboxylate (other name: PB-22);

(8-quinoliny)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);

(8-quinoliny)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);

N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);

N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name: AB-FUBINACA);

1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);

N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: ADB-PINACA);

N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name: AB-CHMINACA);

N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:

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

ENGROSSED

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