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**HOUSE BILL NO. 1564**

Offered January 14, 2015

Prefiled January 7, 2015

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

Patron—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:

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;  
Etoxeridine;  
Furethidine;  
Hydroxypethidine;  
Ketobemidone;  
Levomoramide;  
Levophenacetylmorphan;  
Morpheridine;  
Noracymethadol;  
Norlevorphanol;  
Normethadone;  
Norpipanone;  
Phenadoxone;  
Phenampromide;  
Phenomorphane;  
Phenoperidine;  
Piritramide;  
Proheptazine;  
Properidine;  
Propiram;  
Racemoramide;

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59 Tilidine;

60 Trimeperidine.

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

64 Acetorphine;

65 Acetyldihydrocodeine;

66 Benzylmorphine;

67 Codeine methylbromide;

68 Codeine-N-Oxide;

69 Cyprenorphine;

70 Desomorphine;

71 Dihydromorphine;

72 Drotebanol;

73 Etorphine;

74 Heroin;

75 Hydromorphanol;

76 Methyldesorphine;

77 Methyldihydromorphine;

78 Morphine methylbromide;

79 Morphine methylsulfonate;

80 Morphine-N-Oxide;

81 Myrophine;

82 Nicocodeine;

83 Nicomorphine;

84 Normorphine;

85 Pholcodine;

86 Thebacon.

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

92 Alpha-ethyltryptamine (some trade or other names: Monase;a-ethyl-1H-indole-3-ethanamine;  
93 3-2-aminobutyl] indole; a-ET; AET);

94 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:  
95 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);

96 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:  
97 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);

98 3,4-methylenedioxy amphetamine;

99 5-methoxy-3,4-methylenedioxy amphetamine;

100 3,4,5-trimethoxy amphetamine;

101 Alpha-methyltryptamine (other name: AMT);

102 Bufotenine;

103 Diethyltryptamine;

104 Dimethyltryptamine;

105 4-methyl-2,5-dimethoxyamphetamine;

106 2,5-dimethoxy-4-ethylamphetamine (DOET);

107 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

108 Ibogaine;

109 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);

110 Lysergic acid diethylamide;

111 Mescaline;

112 Parahexyl (some trade or other names: 3-Hexyl-1-hydroxy-7, 8, 9, 10-tetrahydro-6, 6,  
113 9-trimethyl-6H-dibenzo -b,d] pyran; Synhexyl);

114 Peyote;

115 N-ethyl-3-piperidyl benzilate;

116 N-methyl-3-piperidyl benzilate;

117 Psilocybin;

118 Psilocyn;

119 Salvinorin A;

120 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated

121 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;  
 122 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish);  
 123 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;  
 124 2,5-DMA);  
 125 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts  
 126 and salts of isomers;  
 127 3,4-methylenedioxymethamphetamine (also known as N-ethyl-alpha-methyl-3,4  
 128 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);  
 129 N-hydroxy-3,4-methylenedioxymethamphetamine (some other names:  
 130 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);  
 131 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:  
 132 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);  
 133 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;  
 134 paramethoxyamphetamine; PMA);  
 135 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,  
 136 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);  
 137 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy,  
 138 PHP);  
 139 Thiophene analog of phencyclidine (some other names: 1-(2-thienyl) -cyclohexyl]-piperidine,  
 140 2-thienyl analog of phencyclidine, TPCP, TCP);  
 141 1-(2-thienyl)cyclohexylpyrrolidine (other name: TCPy);  
 142 3,4-methylenedioxypyrovalerone (other name: MDPV);  
 143 4-methylmethcathinone (other names: mephedrone, 4-MMC);  
 144 3,4-methylenedioxymethcathinone (other name: methylone);  
 145 Naphthylpyrovalerone (other name: naphyrone);  
 146 4-fluoromethcathinone (other name: flephedrone, 4-FMC);  
 147 4-methoxymethcathinone (other names: methedrone; bk-PMMA);  
 148 Ethcathinone (other name: N-ethylcathinone);  
 149 3,4-methylenedioxyethcathinone (other name: ethylone);  
 150 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone);  
 151 N,N-dimethylcathinone (other name: metamfepramone);  
 152 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);  
 153 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);  
 154 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);  
 155 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);  
 156 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);  
 157 3-fluoromethcathinone (other name: 3-FMC)  
 158 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);  
 159 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);  
 160 4-Methylethcathinone (other name: 4-MEC);  
 161 4-Ethylmethcathinone (other name: 4-EMC);  
 162 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);  
 163 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP);  
 164 Alpha-methylamino-butyrophenone (other name: Buphedrone);  
 165 Alpha-methylamino-valerophenone (other name: Pentedrone);  
 166 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);  
 167 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);  
 168 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,  
 169 25I-NBOMe);  
 170 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);  
 171 4-Fluoromethamphetamine (other name: 4-FMA);  
 172 4-Fluoroamphetamine (other name: 4-FA);  
 173 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);  
 174 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);  
 175 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);  
 176 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);  
 177 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);  
 178 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);  
 179 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);  
 180 (2-aminopropyl)benzofuran (other name: APB);  
 181 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);

- 182 4-chloro-2,5-dimethoxy-N-[2-methoxyphenyl]methyl]-benzeneethanamine (other names:  
 183 2C-C-NBOMe, 25C-NBOMe);  
 184 4-bromo-2,5-dimethoxy-N-[2-methoxyphenyl]methyl]-benzeneethanamine (other names:  
 185 2C-B-NBOMe, 25B-NBOMe);  
 186 Acetoxymethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);  
 187 Benocyclidine (other names: BCP, BTCP);  
 188 Alpha-pyrrolidinobutylphenone (other name: alpha-PBP);  
 189 3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA).  
 190 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture  
 191 or preparation which contains any quantity of the following substances having a depressant effect on the  
 192 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such  
 193 salts, isomers and salts of isomers is possible within the specific chemical designation:  
 194 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;  
 195 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);  
 196 Mecloqualone;  
 197 Methaqualone.  
 198 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture  
 199 or preparation which contains any quantity of the following substances having a stimulant effect on the  
 200 central nervous system, including its salts, isomers and salts of isomers:  
 201 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,  
 202 5-dihydro-5-phenyl-2-oxazolamine);  
 203 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);  
 204 Fenethylline;  
 205 Ethylamphetamine;  
 206 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,  
 207 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;  
 208 Methcathinone (some other names: 2-(methylamino)-propionophenone; ~~alpha-(methylamino)-~~  
 209 ~~propionophenone~~ *alpha-(methylamino)-propionophenone*; 2-(methylamino)-1-phenylpropan-1-one;  
 210 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;  
 211 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);  
 212 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);  
 213 N,N-dimethylamphetamine (other names: N,N-alpha-trimethyl-benzeneethanamine,  
 214 N,N-alpha-trimethylphenethylamine).  
 215 6. Any material, compound, mixture or preparation containing any quantity of the following  
 216 substances:  
 217 N-3-methyl-1-(2-phenethyl)-4-piperidyl]-N-phenylpropanamide (other name: 3-methylfentanyl), its  
 218 optical and geometric isomers, salts, and salts of isomers;  
 219 1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP), its optical isomers, salts and salts of  
 220 isomers;  
 221 1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP), its optical isomers, salts and  
 222 salts of isomers;  
 223 N-1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propionanilide (other names:  
 224 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine), alpha-methylfentanyl);  
 225 N-1-(1-methyl-2-phenethyl)-4-piperidyl]-N-phenylacetamide (other name:  
 226 acetyl-alpha-methylfentanyl), its optical isomers, salts and salts of isomers;  
 227 N-1-(1-methyl-2-2-thienyl)ethyl-4 piperidyl]-N-phenylpropanamide (other name:  
 228 alpha-methylthiofentanyl), its optical isomers, salts and salts of isomers;  
 229 N-1-benzyl-4-piperidyl]N-phenylpropanamide (other name: benzylfentanyl), its optical isomers, salts  
 230 and salts of isomers;  
 231 N-1-(2-hydroxy-2-phenyl) ethyl-4-piperidyl]-N-phenylpropanamide (other name:  
 232 beta-hydroxyfentanyl), its optical isomers, salts and salts of isomers;  
 233 ~~N-3-methyl-1-(2-hydroxy-2-phenethyl)4-piperidyl]N-phenylpropanamide~~  
 234 *N-3-methyl-1-(2-hydroxy-2-phenethyl)4-piperidyl]-N-phenylpropanamide* (other name:  
 235 ~~beta-hydroxy-3-methylfentanyl~~ *beta-hydroxy-3-methylfentanyl*), its optical and geometric isomers, salts and  
 236 salts of isomers;  
 237 N-(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (other name:  
 238 3-methylthiofentanyl), its optical and geometric isomers, salts and salts of isomers;  
 239 N-1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (other name: ~~thienylfentanyl~~ *thienylfentanyl*),  
 240 its optical isomers, salts and salts of isomers;  
 241 N-phenyl-N-1-(2-thienyl)ethyl-4-piperidinyl]-propanamide (other name: thiofentanyl), its optical  
 242 isomers, salts and salts of isomers;  
 243 ~~N-(4-fluorophenyl)-N-1-(2-phenethyl)-4-piperidinyl]-propanamide~~

244 *N*-(4-fluorophenyl)-*N*-1-(2-phenethyl)-4-piperidinyl]-propanamide (other name: para-fluorofentanyl), its  
 245 optical isomers, salts and salts of isomers.

246 7. Any substance that contains one or more cannabimimetic agents or that contains their salts,  
 247 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is  
 248 possible within the specific chemical designation, and any preparation, mixture, or substance containing,  
 249 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.

250 a. "Cannabimimetic agents" includes any substance that is within any of the following structural  
 251 classes:

252 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or  
 253 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;

254 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of  
 255 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not  
 256 substituted on the naphthoyl or naphthyl ring to any extent;

257 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not  
 258 further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to  
 259 any extent;

260 1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not  
 261 further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to  
 262 any extent;

263 3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring,  
 264 whether or not further substituted in the indole ring to any extent, whether or not substituted on the  
 265 phenyl ring to any extent;

266 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring, whether or not further  
 267 substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any  
 268 extent;

269 3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further  
 270 substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any  
 271 extent;

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

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

278 b. The term "cannabimimetic agents" includes:

279 5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);

280 5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);

281 5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);

282 5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);

283 1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);

284 1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);

285 1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);

286 1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);

287 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);

288 (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chr-  
 289 omen-1-ol (other name: HU-210);

290 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);

291 1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);

292 1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);

293 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);

294 1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);

295 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);

296 1-((*N*-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);

297 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);

298 1-[(*N*-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);

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

301 1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);

302 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);

303 1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);

304 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: XLR-11);

305 N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);  
306 N-adamantyl-1-pentylindazole-3-carboxamide (other name: AKB48);  
307 1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);  
308 (8-quinoliny)(1-pentylindol-3-yl)carboxylate (other name: PB-22);  
309 (8-quinoliny)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);  
310 (8-quinoliny)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);  
311 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);  
312 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name:  
313 AB-FUBINACA);  
314 1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (*other name: THJ-2201*);  
315 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name:  
316 ADB-PINACA);  
317 *N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name:*  
318 *AB-CHMINACA)*;  
319 *N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:*  
320 *5-fluoro-AMB)*.  
321 **2. That the provisions of this act may result in a net increase in periods of imprisonment or**  
322 **commitment. Pursuant to § 30-19.1:4, the estimated amount of the necessary appropriation is \$0**  
323 **for periods of imprisonment in state adult correctional facilities and cannot be determined for**  
324 **periods of commitment to the custody of the Department of Juvenile Justice.**