

2015 SESSION

LEGISLATION NOT PREPARED BY DLS
INTRODUCED

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SB1380

15104005D

SENATE BILL NO. 1380

Offered January 20, 2015

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

Patron—Obenshain

Referred to Committee on Education and Health

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;

Etoxidine;

Furethidine;

Hydroxypethidine;

Ketobemidone;

Levomoramide;

Levophenacymorphan;

Morpheridine;

Noracymethadol;

Norlevorphanol;

Normethadone;

Norpipanone;

Phenadoxone;

Phenampromide;

Phenomorphane;

Phenoperidine;

Piritramide;

Proheptazine;

Propiridine;

Propiram;

Racemoramide;

Tilidine;

59 Trimeperidine.

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

63 Acetorphine;

64 Acetyldihydrocodeine;

65 Benzylmorphine;

66 Codeine methylbromide;

67 Codeine-N-Oxide;

68 Cyprenorphine;

69 Desomorphine;

70 Dihydromorphine;

71 Drotebanol;

72 Etorphine;

73 Heroin;

74 Hydromorphanol;

75 Methyldesorphine;

76 Methyldihydromorphine;

77 Morphine methylbromide;

78 Morphine methylsulfonate;

79 Morphine-N-Oxide;

80 Myrophine;

81 Nicocodeine;

82 Nicomorphine;

83 Normorphine;

84 Pholcodine;

85 Thebacon.

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

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

93 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:
94 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);

95 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:
96 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);

97 3,4-methylenedioxy amphetamine;

98 5-methoxy-3,4-methylenedioxy amphetamine;

99 3,4,5-trimethoxy amphetamine;

100 Alpha-methyltryptamine (other name: AMT);

101 Bufotenine;

102 Diethyltryptamine;

103 Dimethyltryptamine;

104 4-methyl-2,5-dimethoxyamphetamine;

105 2,5-dimethoxy-4-ethylamphetamine (DOET);

106 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

107 Ibogaine;

108 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);

109 Lysergic acid diethylamide;

110 Mescaline;

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

113 Peyote;

114 N-ethyl-3-piperidyl benzilate;

115 N-methyl-3-piperidyl benzilate;

116 Psilocybin;

117 Psilocyn;

118 Salvinorin A;

119 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated
120 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;

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

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

optical isomers, salts and salts of isomers.

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

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

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

3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane 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 naphthoyl or naphthyl ring to any extent;

3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not 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-cyclopropylindole 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);

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

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