

# **WEST VIRGINIA LEGISLATURE**

## **2021 REGULAR SESSION**

**Introduced**

### **House Bill 3113**

BY DELEGATES YOUNG, FLUHARTY, PUSHKIN AND

HORNBUCKLE

[Introduced March 12, 2021; Referred to the  
Committee on Health and Human Resources then the  
Judiciary]

1 A BILL to amend and reenact §60A-2-204 of the Code of West Virginia, 1931, as amended,  
 2 relating to removing certain substances from schedule I of the Uniform Controlled  
 3 Substances Act, including marihuana, Psilocybin, and Tetrahydrocannabinols.

*Be it enacted by the Legislature of West Virginia:*

**ARTICLE 2. STANDARDS AND SCHEDULES.**

**§60A-2-204. Schedule I.**

1 (a) Schedule I shall consist of the drugs and other substances, by whatever official name,  
 2 common or usual name, chemical name, or brand name designated, listed in this section including  
 3 their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the  
 4 existence of such isomers, esters, ethers, and salts is possible within the specific chemical  
 5 designation.

6 (b) Opiates.

7 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl) -4-piperidiny]—  
 8 phenylacetamide);

9 Acetylmethadol;

10 Allylprodine;

11 Alphacetylmethadol (except levoalphacetylmethadol also known as levo-alpha-  
 12 acetylmethadol, levomethadyl acetate, or LAAM);

13 Alphameprodine;

14 Alphamethadol;

15 Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propionanilide;  
 16 1-(1-methyl-2-phenylethyl)-4-(( propanilido) piperidine);

17 Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl) ethyl- 4-piperidiny]—  
 18 phenylpropanamide);

19 Benzethidine;

20 Betacetylmethadol;

- 21 Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl) -4- piperidiny]-N-  
22 phenylpropanamide);
- 23 Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2- hydroxy-2-phenethyl)-3-methyl-4-  
24 piperidiny]-N-phenylpropanamide);
- 25 Betameprodine;
- 26 Betamethadol;
- 27 Betaprodine;
- 28 Clonitazene;
- 29 Dextromoramide;
- 30 Diampromide;
- 31 Diethylthiambutene;
- 32 Difenoxin;
- 33 Dimenoxadol;
- 34 Dimepheptanol;
- 35 Dimethylthiambutene;
- 36 Dioxaphetyl butyrate;
- 37 Dipipanone;
- 38 Ethylmethylthiambutene;
- 39 Etonitazene;
- 40 Etoxidine;
- 41 Furethidine;
- 42 Hydroxypethidine;
- 43 Ketobemidone;
- 44 Levomoramide;
- 45 Levophenacymorphan;
- 46 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4- piperidy]-N-phenylpropanamide);

- 47 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4- piperidiny]—phenylpropanamide);
- 48 Morpheridine;
- 49 MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- 50 Noracymethadol;
- 51 Norlevorphanol;
- 52 Normethadone;
- 53 Norpipanone;
- 54 Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2- phenethyl)-4-piperidiny] propanamide);
- 55 PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 56 Phenadoxone;
- 57 Phenampromide;
- 58 Phenomorphan;
- 59 Phenoperidine;
- 60 Piritramide;
- 61 Proheptazine;
- 62 Properidine;
- 63 Propiram;
- 64 Racemoramide;
- 65 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4- piperidiny]-propanamide);
- 66 Tilidine;
- 67 Trimeperidine.
- 68 (c) Opium derivatives:
- 69 Acetorphine;
- 70 Acetyldihydrocodeine;
- 71 Benzylmorphine;
- 72 Codeine methylbromide;

- 73 Codeine-N-Oxide;
- 74 Cyprenorphine;
- 75 Desomorphine;
- 76 Dihydromorphine;
- 77 Drotebanol;
- 78 Etorphine (except HCl Salt);
- 79 Heroin;
- 80 Hydromorphinol;
- 81 Methyldesorphine;
- 82 Methyldihydromorphine;
- 83 Morphine methylbromide;
- 84 Morphine methylsulfonate;
- 85 Morphine-N-Oxide;
- 86 Myrophine;
- 87 Nicocodeine;
- 88 Nicomorphine;
- 89 Normorphine;
- 90 Pholcodine;
- 91 Thebacon.
- 92 (d) Hallucinogenic substances.
- 93 Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-
- 94 indole-3-ethanamine; 3-(2- aminobutyl) indole; alpha-ET; and AET;
- 95 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
- 96 dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;
- 97 4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-
- 98 dimethoxyphenyl)-1-aminoethane; alpha- desmethyl DOB; 2C-B, Nexus;

99 N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the  
100 acronym 25B-NBOMe.

101 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe)

102 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)

103 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-  
104 methylphenethylamine; 2,5-DMA;

105 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

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

107 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-  
108 methylphenethylamine; paramethoxyamphetamine; PMA;

109 5-methoxy-3, 4-methylenedioxy-amphetamine;

110 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-  
111 dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

112 3,4-methylenedioxy amphetamine;

113 3,4-methylenedioxymethamphetamine (MDMA);

114 3,4-methylenedioxy-N-ethylamphetamine (also known as ( ethyl-alpha-methyl-3,4  
115 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

116 N-hydroxy-3,4-methylenedioxyamphetamine (also known as ( hydroxy-alpha-methyl-3,4  
117 (methylenedioxy) phenethylamine, and ( hydroxy MDA);

118 3,4,5-trimethoxy amphetamine;

119 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);

120 Alpha-methyltryptamine (other name: AMT);

121 Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-  
122 hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-  
123 dimethyltryptamine; mappine;

124 Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;

125 Dimethyltryptamine; some trade or other names: DMT;  
126 5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);  
127 Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-  
128 methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepero [5,4-b] indole; Tabernanthe iboga;  
129 Lysergic acid diethylamide;  
130 ~~Marihuana;~~  
131 Mescaline;  
132 Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,  
133 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;  
134 Peyote; meaning all parts of the plant presently classified botanically as Lophophora  
135 williamsii Lemaire, whether growing or not, the seeds thereof, any extract from any part of such  
136 plant, and every compound, manufacture, salts, immediate derivative, mixture, or preparation of  
137 such plant, its seeds or extracts;  
138 N-ethyl-3-piperidyl benzilate;  
139 N-methyl-3-piperidyl benzilate;  
140 Psilocybin;  
141 Psilocyn;  
142 ~~Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or~~  
143 ~~in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives~~  
144 ~~and their isomers with similar chemical structure and pharmacological activity such as the~~  
145 ~~following:~~  
146 ~~delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;~~  
147 ~~delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;~~  
148 ~~delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;~~  
149 ~~(Since nomenclature of these substances is not internationally standardized, compounds~~  
150 ~~of these structures, regardless of numerical designation of atomic positions covered.)~~

151 Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-  
152 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,  
153 cyclohexamine, PCE;

154 Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-  
155 pyrrolidine, PCPy, PHP;

156 Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-  
157 cyclohexyl]-piperidine, 2-thienyl analog of phencyclidine; TCP, TCP;

158 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy.

159 4-methylmethcathinone (Mephedrone);

160 3,4-methylenedioxypropylvalerone (MDPV);

161 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

162 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D)

163 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C)

164 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I)

165 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2)

166 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4)

167 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H)

168 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N)

169 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P)

170 3,4-Methylenedioxy-N-methylcathinone (Methylone)

171 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7, its optical isomers, salts and  
172 salts of isomers)

173 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-  
174 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT)

175 Alpha-methyltryptamine (other name: AMT)

176 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT)

177 Synthetic Cannabinoids as follows:

178 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol {also known as CP  
179 47,497 and homologues};

180 rel-2-[(1S,3R)-3-hydroxycyclohexyl]-5-(2-methylnonan-2-yl)phenol {also known as CP  
181 47,497-C8 homolog};

182 [(6aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-  
183 tetrahydrobenzo[c]chromen-1-ol] {also known as HU-210};  
184 (dexanabinol);

185 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-  
186 tetrahydrobenzo[c]chromen-1-ol {also known as HU-211};

187 1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};

188 1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};

189 (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-015};

190 (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};

191 [1-[2-(4-morpholinyl) ethyl]-1H-indol-3-yl]-1-naphthalenyl-methanone {also known as  
192 JWH-200};

193 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};

194 2-((1S,2S,5S)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl)-5-(2-methyloctan-2-yl)phenol  
195 {also known as CP 55,940};

196 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-122};

197 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-398};

198 (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};

199 1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl) ethanone {also known as  
200 RCS-8};

201 1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);

202 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and

- 203 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).
- 204 Synthetic cannabinoids:
- 205 CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
- 206 YL)phenol);
- 207 HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,
- 208 10A-tetrahydrobenzo[C] chromen-1-OL)];
- 209 HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-
- 210 YL)-6A,7,10,10atetrahydrobenzo[ C]chromen-1-OL);
- 211 JWH-018, 1-pentyl-3-(1-naphthoyl)indole;
- 212 JWH-019, 1-hexyl-3-(1-naphthoyl)indole;
- 213 JWH-073, 1-butyl-3-(1-naphthoyl)indole;
- 214 JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)- Naphthalen-1-ylmethanone;
- 215 JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.]
- 216 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-
- 217 ADB);
- 218 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB);
- 219 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-
- 220 AMB);
- 221 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA);
- 222 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
- 223 (ADB-FUBINACA);
- 224 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
- 225 (MDMB-CHMICA);
- 226 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate
- 227 (MDMB-FUBINACA);
- 228 Tetrahydrocannabinols:

- 229 ~~DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.~~
- 230 ~~DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.~~
- 231 ~~DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.~~
- 232 Synthetic Phenethylamines
- 233 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
- 234 NBOMe);
- 235 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-
- 236 NBOMe);
- 237 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
- 238 NBOMe);
- 239 Synthetic Opioids (including their isomers, esters, ethers, salts and salts of isomers, esters
- 240 and ethers):
- 241 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);
- 242 furanyl fentanyl;
- 243 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
- 244 47700);
- 245 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
- 246 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);
- 247 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also known
- 248 as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-
- 249 hydroxythiofentanyl).
- 250 N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl)
- 251 N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl)
- 252 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl)
- 253 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide
- 254 (also known as U-48800)

255 Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as  
256 U-49900)

257 Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also  
258 known as U-51754)

259 Opioid Receptor Agonist

260 AH-7921 (3,4-dichloro-N-(1dimethylamino)cyclohexylmethyl]benzamide).

261 Naphthoylindoles or any compound containing a 3-(1-Naphthoyl) indole structure with  
262 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole  
263 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall  
264 include the following:

265 JWH 015;

266 JWH 018;

267 JWH 019;

268 JWH 073;

269 JWH 081;

270 JWH 122;

271 JWH 200;

272 JWH 210;

273 JWH 398;

274 AM 2201;

275 WIN 55,212.

276 Naphylmethylindoles or any compound containing a 1indol-3-yl-(1-naphthyl) methane  
277 structure with a substitution at the nitrogen atom of the indole ring whether or not further  
278 substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to  
279 any extent. This shall include, but not be limited to, JWH 175 and JWH 184.

280 Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure with  
281 substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the pyrrole  
282 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall  
283 include, but not be limited to, JWH 147 and JWH 307.

284 Naphthylmethylenes or any compound containing a Naphthylideneindene structure  
285 with substitution at the 3- Position of the indene ring whether or not further substituted in the  
286 indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This  
287 shall include, but not be limited to, JWH 176.

288 Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure with  
289 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole  
290 ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include  
291 the following:

292 RCS-8, SR-18 OR BTM-8;

293 JWH 250;

294 JWH 203;

295 JWH 251;

296 JWH 302.

297 Cyclohexylphenols or any compound containing a 2-(3- hydroxycyclohexyl) phenol  
298 structure with a substitution at the 5-position of the phenolic ring whether or not substituted in the  
299 cyclohexyl ring to any extent. This shall include the following:

300 CP 47,497 and its homologues and analogs;

301 Cannabicyclohexanol;

302 CP 55,940.

303 Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with  
304 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole

305 ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include  
306 the following:

307 AM 694;

308 Pravadoline WIN 48,098;

309 RCS 4;

310 AM 679.

311 [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-  
312 naphthalenymethanone. This shall include WIN 55,212-2.

313 Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol  
314 structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051,  
315 and JWH 133.

316 Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with  
317 substitution at the nitrogen atom of the indole ring whether or not further substituted in the  
318 adamantoyl ring system to any extent. This shall include AM1248.

319 Tetramethylcyclopropylindoles or any compound containing A 3-  
320 tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring  
321 whether or not further substituted in the indole ring to any extent and whether or not substituted  
322 in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.

323 N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.

324 Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as  
325 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV,  
326 and V, not federal Food and Drug Administration approved drug or used within legitimate,  
327 approved medical research. Since nomenclature of these substances is not internationally  
328 standardized, any immediate precursor or immediate derivative of these substances shall be  
329 covered.

330 Tryptamines:

- 331 5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)
- 332 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT)
- 333 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)
- 334 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)
- 335 4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT)
- 336 5-methoxy- $\alpha$ -methyltryptamine (5-MeO-AMT)
- 337 4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT)
- 338 4-hydroxy Diethyltryptamine (4-HO-DET)
- 339 5- methoxy- N,N- diallyltryptamine (5-MeO-DALT)
- 340 4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT)
- 341 4-hydroxy Diethyltryptamine (4-HO-DET)
- 342 (e) Depressants.
- 343 Mecloqualone;
- 344 Methaqualone.
- 345 (f) Stimulants.
- 346 Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
- 347 dihydro-5-phenyl-2-oxazolamine;
- 348 Cathinone; some trade or other names: 2-amino-1-phenyl-1- propanone, alpha-
- 349 aminopropiophenone, 2-aminopropiophenone, and norephedrone;
- 350 Fenethylline;
- 351 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
- 352 isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-
- 353 (methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—
- 354 methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxyprovalerone and/or
- 355 mephedrone;3,4-methylenedioxyprovalerone (MPVD); ephedrone; N-methylcathinone;
- 356 methylcathinone; AL-464; AL-422; AL- 463 and UR1432;

- 357 (+-) cis-4-methylaminorex; ((+)-)cis-4,5-dihydro-4-methyl- 5-phenyl-2-oxazolamine);
- 358 N-ethylamphetamine;
- 359 N,N-dimethylamphetamine; also known as N,N-alpha- trimethyl-benzeneethanamine;
- 360 N,N-alpha-trimethylphenethylamine.
- 361 Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts, and
- 362 salts of isomers.
- 363 Substituted amphetamines:
- 364 2-Fluoroamphetamine
- 365 3-Fluoroamphetamine
- 366 4-Fluoroamphetamine
- 367 2-chloroamphetamine
- 368 3-chloroamphetamine
- 369 4-chloroamphetamine
- 370 2-Fluoromethamphetamine
- 371 3-Fluoromethamphetamine
- 372 4-Fluoromethamphetamine
- 373 4-chloromethamphetamine
- 374 (g) Temporary listing of substances subject to emergency scheduling. Any material,
- 375 compound, mixture, or preparation which contains any quantity of the following substances:
- 376 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts,
- 377 and salts of isomers.
- 378 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
- 379 isomers, salts, and salts of isomers.
- 380 N-benzylpiperazine, also known as BZP.
- 381 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);

382 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-  
383 butyramide);

384 Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);

385 Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-  
386 acetamide);

387 3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-  
388 phenylbutyramide);

389 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-  
390 yl)butyramide);

391 Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);

392 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-  
393 carboxamide);

394 Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).

395 (h) The following controlled substances are included in Schedule I:

396 Synthetic Cathinones or any compound, except bupropion or compounds listed under a  
397 different schedule, or compounds used within legitimate and approved medical research,  
398 structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic  
399 or fused polycyclic ring systems, whether or not the compound is further modified in any of the  
400 following ways:

401 By substitution in the ring system to any extent with Alkyl, alkylendioxy, alkoxy, haloalkyl,  
402 hydroxyl, or halide Substituents whether or not further substituted in the ring system by one or  
403 more other univalent substituents.

404 By substitution at the 3-position with an acyclic alkyl substituent.

405 By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl or methoxybenzyl  
406 groups.

407 By inclusion of the 2-amino nitrogen atom in a cyclic structure.

408           Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as  
409 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV,  
410 and V, not federal Food and Drug Administration approved drug or used within legitimate,  
411 approved medical research.

NOTE: The purpose of this bill is to remove certain substances from schedule I of the Uniform Controlled Substances Act, including marihuana, Psilocybin, and Tetrahydrocannabinols.

Strike-throughs indicate language that would be stricken from a heading or the present law and underscoring indicates new language that would be added.