

WEST VIRGINIA LEGISLATURE

2026 REGULAR SESSION

Introduced

House Bill 5146

By Delegate Young

[Introduced February 03, 2026; 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-piperidinyl]-
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-piperidinyl]-
18 phenylpropanamide);

19 Benzethidine;

20 Betacetylmethadol;

21 Beta-hydroxyfentanyl(N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-
22 phenylpropanamide);
23 Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-
24 piperidinyl]-N-phenylpropanamide);
25 Betameprodine;
26 Betamethadol;
27 Betaprodine;
28 Brorphine (1-(1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2H-benzo[d]imidazol-
29 2-one);
30 Clonitazene;
31 Dextromoramide;
32 Diamprodine;
33 Diethylthiambutene;
34 Difenoxyin;
35 Dimenoxadol;
36 Dimepheptanol;
37 Dimethylthiambutene;
38 Dioxaphetyl butyrate;
39 Dipipanone;
40 Ethylmethylthiambutene;
41 Etonitazene;
42 Etoxeridine;
43 Fentanyl analog or derivative, as that term is defined in article one of this chapter:
44 *Provided*, That fentanyl and carfentanil remains a Schedule II substance, as set forth in W. Va.
45 Code §60A-2-206;
46 Furethidine;

47 Hydroxypethidine;
48 Ketobemidone;
49 Levomoramide;
50 Levophenacylmorphan;
51 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);
52 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4-piperidinyl]-phenylpropanamide);
53 Morpheridine;
54 N-Methylnorfentanyl (N-(1-Methyl-4-piperidinyl)-N-phenyl-propanamide,
55 monohydrochloride);
56
57 MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
58 Noracymethadol;
59 Norlevorphanol;
60 Normethadone;
61 Norpipanone;
62 Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide);
63 PEPAP(1-(2-phenethyl)-4-phenyl-4-acetoxytropine);
64 Phenadoxone;
65 Phenampromide;
66 Phenomorphan;
67 Phenoperidine;
68 Piritramide;
69 Proheptazine;
70 Properidine;
71 Propiram;
72 Racemoramide;

73 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);
74 Tilidine;
75 Trimeperidine.
76 (c) Opium derivatives,
77 Acetorphine;
78 Acetyldihydrocodeine;
79 Benzylmorphine;
80 Codeine methylbromide;
81 Codeine-N-Oxide;
82 Cyprenorphine;
83 Desomorphine;
84 Dihydromorphine;
85 Drotebanol;
86 Etorphine (except HCl Salt);
87 Heroin;
88 Hydromorphenol;
89 Methyldesorphine;
90 Methyldihydromorphine;
91 Morphine methylbromide;
92 Morphine methylsulfonate;
93 Morphine-N-Oxide;
94 Myrophine;
95 Nicocodeine;
96 Nicomorphine;
97 Normorphine;
98 Pholcodine;

99 Thebacon.

100 (d) Hallucinogenic substances.

101 Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-

102 indole-3-ethanamine; 3-(2- aminobutyl) indole; alpha-ET; and AET;

103 1-(4-methoxyphenyl)-N-methylpropan-2-amine (other names: para-methoxymethamphetamine,

104 PMMA);

105 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-

106 dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;

107 4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-

108 dimethoxyphenyl)-1-aminoethane; alpha- desmethyl DOB; 2C-B, Nexus;

109 N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the

110 acronym 25B-NBOMe;

111 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe);

112 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe);

113 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-

114 methylphenethylamine; 2,5-DMA;

115 2,5-dimethoxy-4-ethylamphetamine; some trade or other names: DOET;

116 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

117 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-

118 methylphenethylamine; paramethoxyamphetamine; PMA;

119 3-Hydroxy-phencyclidine (other name hydroxy PCP);

120 5-methoxy-3, 4-methylenedioxy-amphetamine;

121 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-

122 dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

123 3,4-methylenedioxy amphetamine;

124 3,4-methylenedioxymethamphetamine (MDMA);

125 3,4-methylenedioxy-N-ethylamphetamine (also known as (ethyl-alpha-methyl-3,4
126 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);
127 N-hydroxy-3,4-methylenedioxyamphetamine (also known as (hydroxy-alpha-methyl-3,4
128 (methylenedioxy) phenethylamine, and (hydroxy MDA);
129 3,4,5-trimethoxy amphetamine;
130 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
131 Alpha-methyltryptamine (other name: AMT);
132 Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-
133 hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-
134 dimethyltryptamine; mappine;
135 Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;
136 Dimethyltryptamine; some trade or other names: DMT;
137 5-Methoxy-N,N-disopropyltryptamine (5-MeO-DIPT);
138 Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-
139 methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;
140 Lysergic acid diethylamide;
141 Marihuana; Marijuana (Cannabis, sp.);
142 Mescaline;
143 Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,
144 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;
145 Peyote; meaning all parts of the plant presently classified botanically as Lophophora
146 williamsii Lemaire, whether growing or not, the seeds thereof, any extract from any part of such
147 plant, and every compound, manufacture, salts, immediate derivative, mixture, or preparation of
148 such plant, its seeds or extracts;
149 N-ethyl-3-piperidyl benzilate;
150 N-methyl-3-piperidyl benzilate;

151 Psilocybin;

152 Psilocyn;

153 Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or

154 in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives

155 and their isomers with similar chemical structure and pharmacological activity including, but not

156 limited to the following:

157 -delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;

158 -delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;

159 -delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;

160 -delta-8 Cis or trans tetrahydrocannabinol and its optical isomers; and

161 -delta-10 Cis or trans tetrahydrocannabinol and its optical isomers;

162 (Since nomenclature of these substances is not internationally standardized, compounds

163 of these structures, regardless of numerical designation of atomic positions covered.)

164 -Delta-8-tetrahydrocannabinol-O (delta-8-THC-O), Delta-9-tetrahydrocannabinol (delta-9-

165 THC-O) and Synthetic and non naturally occurring cannabinoids.

166 The provisions of this section related to tetrahydrocannabinols are inapplicable to

167 products or substances lawfully manufactured, distributed, or possessed under the provisions of §

168 19-12E-1 *et seq.* and Chapter 16H of this code.

169 Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-

170 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,

171 cyclohexamine, PCE;

172 Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-

173 pyrrolidine, PCPy, PHP;

174 Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-

175 cyclohexyl]-piperidine, 2-thienyl analog of phencyclidine; TPCP, TCP;

176 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy;

177 4-methylmethcathinone (Mephedrone);
178 3,4-methylenedioxypyrovalerone (MDPV);
179 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);
180 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);
181 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);
182 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);
183 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);
184 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);
185 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);
186 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);
187 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);
188 3,4-Methylenedioxy-N-methylcathinone (Methylone);
189 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7, its optical isomers, salts and
190 salts of isomers;
191 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-
192 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);
193 Alpha-methyltryptamine (other name: AMT);
194 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
195 Synthetic Cannabinoids as follows:
196 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol { also known as CP
197 47,497 and homologues} ;
198 rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl)phenol { also known as CP
199 47,497-C8 homolog} ;
200 [(6aR)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7,10,10a-
201 tetrahydrobenzo[c]chromen-1-ol)] { also known as HU-210} ;
202 (dexanabinol);

203 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
204 tetrahydrobenzol[c]chromen-1-ol) { also known as HU-211} ;
205 1-Pentyl-3-(1-naphthoyl)indole { also known as JWH-018} ;
206 1-Butyl-3-(1-naphthoyl)indole { also known as JWH-073} ;
207 (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone { also known as JWH-015} ;
208 (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone { also known as JWH-019} ;
209 [1-[2-(4-morpholinyl) ethyl] -1H-indol-3-yl]-1-naphthalenyl-methanone { also known as
210 JWH-200} ;
211 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone { also known as JWH-250} ;
212 2-((1S,2S,5S)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl) -5-(2-methyloctan-2-yl)phenol {
213 also known as CP 55,940} ;
214 (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl)-methanone { also known as JWH-
215 122};
216 (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl)-methanone { also known as JWH-398};
217 (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone { also known as RCS-4} ;
218 1-(1-(2-cyclohexylethyl) -1H-indol-3-yl) -2-(2-methoxyphenyl) ethanone { also known as
219 RCS-8} ;
220 1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);
221 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and
222 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).
223 Synthetic cannabinoids:
224 CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
225 YL)phenol);
226 HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,
227 10A-tetrahydrobenzo[C] chromen-1-OL)];
228 HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-

229 YL)-6A,7,10,10atetrahydrobenzo[C]chromen-1-OL);
230 JWH-018, 1-pentyl-3-(1-naphthoyl)indole;
231 JWH-019, 1-hexyl-3-(1-naphthoyl)indole;
232 JWH-073, 1-butyl-3-(1-naphthoyl)indole;
233 JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)- Naphthalen-1-ylmethanone;
234 JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.]
235 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-
236 ADB);
237 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB);
238 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-
239 AMB);
240 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA);
241 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
242 (ADB-FUBINACA);
243 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
244 (MDMB-CHMICA);
245 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (MDMB-
246 FUBINACA);
247 Tetrahydrocannabinols:
248 ~~DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.~~
249 ~~DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.~~
250 ~~DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.~~
251 Synthetic Phenethylamines
252 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
253 NBOMe);
254 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-

255 NBOMe);

256 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-

257 NBOMe);

258 Synthetic Opioids (including their isomers, esters, ethers, salts and salts of isomers, esters

259 and ethers):

260 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);

261 furanyl fentanyl;

262 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-

263 47700);

264 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-

265 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);

266 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also known

267 as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-

268 hydroxythiofentanyl);

269 N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl);

270 N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl);

271 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl);

272 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide

273 (also known as U-48800);

274 Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as

275 U-49900);

276 Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also

277 known as U-51754);

278 2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine

279 (butonitazene);

280 2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine (etodesnitazene);

281 N,N-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine
282 (flunitazene);
283 N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine
284 (metodesnitazene);
285 N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine
286 (metonitaze);
287 2-(4-ethoxybenzyl)5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1 H-benzimidazole (N-pyrrolidino
288 etonitazene, etonitazepyne);
289 N,N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine
290 (protonitazene);
291 N-pyrrolidino etonitazene;
292 Etodesnitazene;
293 Isotonitazene;
294 Protonitazene;
295 Metonitazene;
296 Butonitazene;
297 Metodesnitazene;
298 Flunitazene;
299 Opioid Receptor Agonist
300 2-Methyl AP-237 (1-(2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl)butan-1-one)
301 AH-7921 (3,4-dichloro-N-(1dimethylamino)cyclohexylmethyl]benzamide).
302 Naphthoylindoles or any compound containing a 3-(1-Naphthoyl) indole structure with
303 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole
304 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall
305 include the following:
306 JWH 015;

307 JWH 018;

308 JWH 019;

309 JWH 073;

310 JWH 081;

311 JWH 122;

312 JWH 200;

313 JWH 210;

314 JWH 398;

315 AM 2201; and

316 WIN 55,212.

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

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

325 Naphthylmethylindenes or any compound containing a Naphthylideneindene structure
326 with substitution at the 3- Position of the indene ring whether or not further substituted in the
327 indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This
328 shall include, but not be limited to, JWH 176.

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

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

334 JWH 250;

335 JWH 203;

336 JWH 251; and

337 JWH 302.

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

341 CP 47,497 and its homologues and analogs;

342 Cannabicyclohexanol; and

343 CP 55,940.

344 Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include the following:

348 AM 694;

349 Pravadoline WIN 48,098;

350 RCS 4; and

351 AM 679.

352 [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-naphthalenymethanone. This shall include WIN 55,212-2.

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

357 Adamantoylindoles or any compound containing a 3-(1-Adamantoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the

359 adamantoyl ring system to any extent. This shall include AM1248.

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

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

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

370 Tryptamines:

371 5-methoxy-N-methyl-N-isopropyltryptamine (5-MeO-MiPT);

372 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT);

373 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT);

374 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);

375 4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT);

376 5-methoxy- α -methyltryptamine (5-MeO-AMT);

377 4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT);

378 4-hydroxy Diethyltryptamine (4-HO-DET);

379 5-methoxy-N,N-diallyltryptamine (5-MeO-DALT);

380 4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT);

381 4-hydroxy Diethyltryptamine (4-HO-DET);

382 FDU-PB-22 (1-Naphthyl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate);

383 FUB-PB-22 (Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate);

384 5-Fluoro-MN-24 (1-(5-Fluoropentyl)-N-(naphthalen-1-yl)-1H-indole-3-carboxamide);

385 MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
386 SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
387 SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
388 Methyl-Ethylaminopentiophenone;
389 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
390 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
391 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
392 carboxamide);
393 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
394 methylbutanoat);
395 MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
396 SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
397 SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
398 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
399 Methyl-Ethylaminopentiophenone;
400 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
401 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
402 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
403 carboxamide);
404 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
405 methylbutanoat);
406 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
407 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-
408 a][1,4]benzodiazepine);
409 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
410 benzodiazepin-2-one);

411 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
412 a][1,4]diazepine);

413 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
414 a][1,4]benzodiazepine);

415 Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);

416 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
417 a][1,4]benzodiazepine);

418 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-
419 a][1,4]diazepine);

420 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
421 one);

422 Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); and

423 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
424 a][1,4]benzodiazepine).

425 (e) Depressants.

426 4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-
427 carboxamide);

428 Alpha-Phenylacetoacetonitrile (3-Oxo-2-phenylbutanenitrile);

429 2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-cyclohexanone,
430 monohydrochloride);

431 4-MEAP (2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one);

432 Mecloqualone;

433 Methaqualone;

434 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);

435 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-
436 a][1,4]benzodiazepine);

437 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
438 benzodiazepin-2-one);
439 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
440 a][1,4]diazepine);
441 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
442 a][1,4]benzodiazepine);
443 Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
444 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
445 a][1,4]benzodiazepine);
446 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-
447 a][1,4]diazepine);
448 gamma-hydroxybutyric acid (some other names include GHB; gamma-hydroxybutyrate; 4-
449 hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
450 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
451 one);
452 Nitrazepam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
453 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
454 a][1,4]benzodiazepine);
455 Diclazepam (7-Chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-
456 one); and
457 Deschloroetizolam (2-Ethyl-9-methyl-4-phenyl-6H-thieno[3,2-f][1,2,4]triazolo[4,3-
458 a][1,4]diazepine);
459 (f) Stimulants.
460 Aminorex; some other names: aminoxyphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
461 dihydro-5-phenyl-2-oxazolamine;
462 4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine; 4-

463 methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine);

464 Cathinone; some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-

465 aminopropiophenone, 2-aminopropiophenone and norephedrone;

466 Ethylphenidate (ethyl 2-phenyl-2-(piperidin-2-yl)acetate);

467 Fenethylline;

468 Mesocarb (N-phenyl-N'-(3-(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-yl)carbamimidate);

469 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical

470 isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-

471 (methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-

472 methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or

473 mephedrone; 3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone;

474 methylcathinone; AL-464; AL-422; AL-463 and UR1432;

475 (+-) cis-4-methylaminorex; ((+)-)cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);

476 N-ethylamphetamine;

477 N,N-dimethylamphetamine; also known as N,N-alpha-trimethyl-benzeneethanamine;

478 N,N-alpha-trimethylphenethylamine;

479 Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and

480 salts of isomers;

481 Substituted amphetamines:

482 2-Fluoroamphetamine;

483 3-Fluoroamphetamine;

484 4-Fluoroamphetamine;

485 2-chloroamphetamine;

486 3-chloroamphetamine;

487 4-chloroamphetamine;

488 2-Fluoromethamphetamine;

489 3-Fluoromethamphetamine;
490 4-Fluoromethamphetamine;
491 4-chloromethamphetamine;
492 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
493 Alpha-PHP (1-Phenyl-2-(pyrrolidin-1-yl)hexan-1-one);
494 MPHP (1-(4-Methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);
495 PV8 (1-Phenyl-2-(pyrrolidin-1-yl)heptan-1-one);
496 4-Chloro-Alpha-PVP (1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one);
497 N-Ethylhexedrone (2-(Ethylamino)-1-phenylhexan-1-one);
498 Methoxetamine (2-(Ethylamino)-2-(3-methoxyphenyl)-cyclohexanone); and
499 3-Fluorophenmetrazine (2-(3-Fluorophenyl)-3-methylmorpholine);
500 (g) Temporary listing of substances subject to emergency scheduling. Any material,
501 compound, mixture, or preparation which contains any quantity of the following substances:
502 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts,
503 and salts of isomers;
504 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
505 isomers, salts, and salts of isomers.
506 N-benzylpiperazine, also known as BZP;
507 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);
508 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-
509 butyramide);
510 Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);
511 Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-
512 acetamide);
513 3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-
514 phenylbutyramide);

515 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
516 yl)butyramide);

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

518 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
519 carboxamide); and

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

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

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

527 By substitution in the ring system to any extent with Alkyl, alkylenedioxy, alkoxy, haloalkyl,
528 hydroxyl, or halide Substituents whether or not further substituted in the ring system by one or
529 more other univalent substituents;

530 By substitution at the 3-position with an acyclic alkyl substituent;

531 By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl or methoxybenzyl
532 groups;

533 By inclusion of the 2-amino nitrogen atom in a cyclic structure; or

534 Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as
535 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV, and
536 V, not federal Food and Drug Administration approved drug or used within legitimate, approved
537 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.