

WEST VIRGINIA LEGISLATURE

2020 REGULAR SESSION

Originating

Senate Bill 847

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[Originating in the Committee on the Judiciary;
reported on February 24, 2020]

1 A BILL to amend and reenact §60A-2-204 and §60A-2-212 of the Code of West Virginia, 1931,
2 as amended, all relating to updating the controlled substance lists in Schedules I and V.

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 N-Methylnorfentanyl (N-(1-Methyl-4-piperidiny)-N-phenyl-propanamide,
51 monohydrochloride);
52 Norfentanyl (N-Phenyl-N-4-piperidiny-propanamide);
53 Noracymethadol;
54 Norlevorphanol;
55 Normethadone;
56 Norpipanone;
57 Para-fluorofentanyl (N-[4-fluorophenyl]-N-[1-(2- phenethyl)-4-piperidiny] propanamide);
58 PEPAP(1-[2-phenethyl]-4-phenyl-4-acetoxypiperidine);
59 Phenadoxone;
60 Phenampromide;
61 Phenomorphan;
62 Phenoperidine;
63 Piritramide;
64 Proheptazine;
65 Properidine;
66 Propiram;
67 Racemoramide;
68 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4- piperidiny]-propanamide);
69 Tilidine;
70 Trimeperidine.
71 (c) *Opium derivatives.* —
72 Acetorphine;

- 73 Acetyldihydrocodeine;
- 74 Benzylmorphine;
- 75 Codeine methylbromide;
- 76 Codeine-N-Oxide;
- 77 Cyprenorphine;
- 78 Desomorphine;
- 79 Dihydromorphine;
- 80 Drotebanol;
- 81 Etorphine (except HCl Salt);
- 82 Heroin;
- 83 Hydromorphanol;
- 84 Methyldesorphine;
- 85 Methyldihydromorphine;
- 86 Morphine methylbromide;
- 87 Morphine methylsulfonate;
- 88 Morphine-N-Oxide;
- 89 Myrophine;
- 90 Nicocodeine;
- 91 Nicomorphine;
- 92 Normorphine;
- 93 Pholcodine;
- 94 Thebacon.
- 95 (d) *Hallucinogenic substances.* —
- 96 Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-
- 97 indole-3-ethanamine; 3-(2-aminobutyl) indole; alpha-ET; and AET;

98 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
99 dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;

100 4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-
101 dimethoxyphenyl)-1-aminoethane; alpha-desmethyl DOB; 2C-B, Nexus;

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

104 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe)

105 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)

106 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-
107 methylphenethylamine; 2,5-DMA;

108 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

109 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

110 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-
111 methylphenethylamine; paramethoxyamphetamine; PMA;

112 3-Hydroxy-phencyclidine (other name hydroxy PCP)

113 3-methoxy-phencyclidine (other name: Methoxy PCP);

114 5-methoxy-3, 4-methylenedioxy-amphetamine;

115 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-
116 dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

117 3,4-methylenedioxy amphetamine;

118 3,4-methylenedioxymethamphetamine (MDMA);

119 3,4-methylenedioxy-N-ethylamphetamine (also known as (ethyl-alpha-methyl-3,4
120 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

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

123 7-Hydroxymitragynine;

124 3,4,5-trimethoxy amphetamine;

125 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);

126 Alpha-methyltryptamine (other name: AMT);

127 Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole;

128 3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N- dimethyltryptamine;

129 mappine;

130 Diethyltryptamine; some trade and other names: N, N-Diethyltryptamine; DET;

131 Dimethyltryptamine; some trade or other names: DMT;

132 5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);

133 Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-

134 methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;

135 Lysergic acid diethylamide;

136 Marihuana *Marijuana (Cannabis, sp.)*;

137 Mescaline;

138 Mitragynine;

139 Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,

140 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;

141 Peyote; meaning all parts of the plant presently classified botanically as *Lophophora*

142 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from any part of such

143 plant, and every compound, manufacture, salts, immediate derivative, mixture, or preparation of

144 such plant, its seeds, or extracts;

145 N-ethyl-3-piperidyl benzilate;

146 N-methyl-3-piperidyl benzilate;

147 Psilocybin;

148 Psilocyn;

149 Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or
150 in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives
151 and their isomers with similar chemical structure and pharmacological activity such as the
152 following:

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

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

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

156 (Since nomenclature of these substances is not internationally standardized, compounds
157 of these structures, regardless of numerical designation of atomic positions covered.);

158 Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-
159 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,
160 cyclohexamine, PCE;

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

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

165 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy;

166 4-methylmethcathinone (Mephedrone);

167 3,4-methylenedioxypropylvalerone (MDPV);

168 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

169 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

170 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

171 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

172 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

173 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

174 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

175 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);
176 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);
177 3,4-Methylenedioxy-N-methylcathinone (Methylone);
178 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7, its optical isomers, salts and
179 salts of isomers;
180 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-
181 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);
182 Alpha-methyltryptamine (other name: AMT);
183 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
184 Synthetic Cannabinoids as follows:
185 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol {also known as CP
186 47,497 and homologues};
187 rel-2-[(1S,3R)-3-hydroxycyclohexyl]-5-(2-methylnonan-2-yl)phenol {also known as CP
188 47,497-C8 homolog};
189 [(6aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
190 tetrahydrobenzo[c]chromen-1-ol] {also known as HU-210};
191 (dexanabinol);
192 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
193 tetrahydrobenzo[c]chromen-1-ol {also known as HU-211};
194 1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};
195 1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};
196 (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-015};
197 (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};
198 [1-[2-(4-morpholinyl)ethyl]-1H-indol-3-yl]-1-naphthalenyl-methanone {also known as
199 JWH-200};
200 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};

201 2-((1S,2S,5S)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl)-5-(2-methyloctan-2-yl)phenol
202 {also known as CP 55,940};
203 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-122};
204 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-398};
205 (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};
206 1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl)ethanone {also known as
207 RCS-8};
208 1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);
209 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and
210 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).
211 Synthetic cannabinoids:
212 CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
213 YL)phenol);
214 HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,
215 10A-tetrahydrobenzo[C]chromen-1-OL)];
216 HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-
217 YL)-6A,7,10,10A-tetrahydrobenzo[C]chromen-1-OL);
218 JWH-018, 1-pentyl-3-(1-naphthoyl)indole;
219 JWH-019, 1-hexyl-3-(1-naphthoyl)indole;
220 JWH-073, 1-butyl-3-(1-naphthoyl)indole;
221 JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)-Naphthalen-1-ylmethanone;
222 JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.];
223 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-
224 ADB);
225 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB);

226 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-
227 AMB);
228 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA);
229 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
230 (ADB-FUBINACA);
231 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
232 (MDMB-CHMICA);
233 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate
234 (MDMB-FUBINACA);
235 Tetrahydrocannabinols:
236 DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.
237 DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.
238 DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.
239 Synthetic Phenethylamines:
240 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
241 NBOMe);
242 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-
243 NBOMe);
244 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
245 NBOMe);
246 Synthetic Opioids (~~including~~ including their isomers, esters, ethers, salts, and salts of
247 isomers, esters, and ethers):
248 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);
249 furanyl fentanyl;
250 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
251 47700);

252 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
253 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);

254 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also known
255 as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidiny]-N-phenylpropanamide, (beta-
256 hydroxythiofentanyl);

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

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

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

260 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide
261 (also known as U-48800);

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

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

266 *Opioid Receptor Agonist.* —

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

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

272 JWH 015;

273 JWH 018;

274 JWH 019;

275 JWH 073;

276 JWH 081;

277 JWH 122;

278 JWH 200;

279 JWH 210;

280 JWH 398;

281 AM 2201;

282 WIN 55,212.

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

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

291 Naphthylmethyloindenes or any compound containing a Naphthylideneindene structure
292 with substitution at the 3- Position of the indene ring whether or not further substituted in the
293 indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This
294 shall include, but not be limited to, JWH 176.

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

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

300 JWH 250;

301 JWH 203;

302 JWH 251;

303 JWH 302.

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

307 CP 47,497 and its homologues and analogs;

308 Cannabicyclohexanol;

309 CP 55,940.

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

314 AM 694;

315 Pravadoline WIN 48,098;

316 RCS 4;

317 AM 679.

318 [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-
319 naphthalenymethanone. This shall include WIN 55,212-2.

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

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

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

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

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

337 Tryptamines:

338 5-methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT);

339 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT);

340 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT);

341 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);

342 4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT);

343 5-methoxy- α -methyltryptamine (5-MeO-AMT);

344 4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT);

345 4-hydroxy Diethyltryptamine (4-HO-DET);

346 5-methoxy- N,N- diallyltryptamine (5-MeO-DALT);

347 4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT);

348 ~~4-hydroxy Diethyltryptamine (4-HO-DET)~~

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

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

351 5-Fluoro-MN-24 (1-(5-Fluoropentyl)-N-(naphthalen-1-yl)-1H-indole-3-carboxamide);

352 MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);

353 SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);

354 SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);

355 Methyl-Ethylaminopentiophenone;

356 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
357 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
358 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
359 carboxamide);
360 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
361 methylbutanoat);
362 MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
363 SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
364 SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
365 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
366 Methyl-Ethylaminopentiophenone;
367 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
368 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
369 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-
370 3- carboxamide);
371 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
372 methylbutanoat);
373 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
374 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-
375 a][1,4]benzodiazepine);
376 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
377 benzodiazepin-2-one);
378 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
379 a][1,4]diazepine);
380 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
381 a][1,4]benzodiazepine);

382 Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
383 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
384 a][1,4]benzodiazepine);
385 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzof[1,2,4]triazolo[4,3-
386 a][1,4]diazepine);
387 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
388 one) ;
389 Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
390 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
391 a][1,4]benzodiazepine).
392 (e) *Depressants.* —
393 4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-2- yl)-1H-indazole-3-
394 carboxamide);
395 Alpha-Phenylacetoacetonitrile (3-Oxo-2-phenylbutanenitrile);
396 2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-cyclohexanone,
397 monohydrochloride);
398 4-MEAP (2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one);
399 Mecloqualone;
400 Methaqualone;
401 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
402 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3
403 a][1,4]benzodiazepine);
404 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
405 benzodiazepin-2-one);
406 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
407 a][1,4]diazepine);

408 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
409 a][1,4]benzodiazepine);
410 Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
411 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
412 a][1,4]benzodiazepine);
413 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzof[1,2,4]triazolo[4,3-
414 a][1,4]diazepine);
415 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
416 one);
417 Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
418 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
419 a][1,4]benzodiazepine);
420 Declazepam (7-Chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2H-
421 1,4-benzodiazepin-2-one);
422 Deschloroetizolam (2-Ethyl-9-methyl-4-phenyl-6H-thieno[3,2- f][1,2,4]triazolo[4,3-
423 a][1,4]diazepine);
424 (f) *Stimulants.* —
425 Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
426 dihydro-5-phenyl-2-oxazolamine;
427 Cathinone; some trade or other names: 2-amino-1-phenyl-1- propanone, alpha-
428 aminopropiophenone, 2-aminopropiophenone and norephedrone;
429 Fenethylline;
430 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
431 isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-
432 (methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—
433 methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxyprovalerone and/or

434 mephedrone;3,4-methylenedioxypropylvalerone (MPVD); ephedrone; N-methylcathinone;
435 methylcathinone; AL-464; AL-422; AL- 463 and UR1432;
436 (+-) cis-4-methylaminorex; ((+)-)cis-4,5-dihydro-4-methyl- 5-phenyl-2-oxazolamine);
437 N-ethylamphetamine;
438 N,N-dimethylamphetamine; also known as N,N-alpha- trimethyl-benzeneethanamine;
439 N,N-alpha-trimethylphenethylamine.
440 Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and
441 salts of isomers.
442 Substituted amphetamines:
443 2-Fluoroamphetamine;
444 3-Fluoroamphetamine;
445 4-Fluoroamphetamine;
446 2-chloroamphetamine;
447 3-chloroamphetamine;
448 4-chloroamphetamine;
449 2-Fluoromethamphetamine;
450 3-Fluoromethamphetamine;
451 4-Fluoromethamphetamine;
452 4-chloromethamphetamine;
453 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
454 Alpha-PHP (1-Phenyl-2-(pyrrolidin-1-yl)hexan-1-one);
455 MPHP (1-(4-Methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);
456 PV8 (1-Phenyl-2-(pyrrolidin-1-yl)heptan-1-one);
457 4-Chloro-Alpha-PVP (1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one);
458 N-Ethylhexedrone (2-(Ethylamino)-1-phenylhexan-1-one);
459 Methoxetamine (2-(Ethylamino)-2-(3-methoxyphenyl)-cyclohexanone);

460 3-Fluorophenmetrazine (2-(3-Fluorophenyl)-3-methylmorpholine);
461 (g) Temporary listing of substances subject to emergency scheduling. Any material,
462 compound, mixture, or preparation which contains any quantity of the following substances:
463 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts,
464 and salts of isomers;
465 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
466 isomers, salts and salts of isomers;
467 N-benzylpiperazine, also known as BZP;
468 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);
469 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-
470 butyramide);
471 Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);
472 Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-
473 acetamide);
474 3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-
475 phenylbutyramide);
476 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
477 yl)butyramide);
478 Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);
479 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
480 carboxamide);
481 Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).
482 (h) The following controlled substances are included in Schedule I:
483 Synthetic cathinones or any compound, except bupropion or compounds listed under a
484 different schedule, or compounds used within legitimate and approved medical research,
485 structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic

486 or fused polycyclic ring systems, whether or not the compound is further modified in any of the
487 following ways:

488 By substitution in the ring system to any extent with alkyl, alkylendioxy, alkoxy, haloalkyl,
489 hydroxyl or halide substituents whether or not further substituted in the ring system by one or
490 more other univalent substituents.

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

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

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

495 Any other synthetic chemical compound that is a cannabinoid receptor type 1 agonist as
496 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV,
497 and V, not a federal Food and Drug Administration-approved drug, or used within legitimate,
498 approved medical research.

§60A-2-212. Schedule V.

1 (a) Schedule V 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. Unless
3 specifically excepted or unless listed in another schedule, any material, compound, mixture, or
4 preparation which contains any quantity of the following substances, including their isomers,
5 esters, ethers, salts, and salts of isomers, esters and ethers, whenever the existence of such
6 isomers, esters, ethers, and salts is possible within the specific chemical designation.

7 (b) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound,
8 mixture or preparation containing any of the following narcotic drugs or their salts calculated as
9 the free anhydrous base or alkaloid in limited quantities as set forth below, which shall include
10 one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the
11 compound, mixture, or preparation valuable medicinal qualities other than those possessed by
12 the narcotic drug alone.

13 Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;
14 Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams;
15 Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams;
16 Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine
17 sulfate per dosage unit;
18 Not more than 100 milligrams of opium per 100 milliliters or per 100 grams;
19 Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine
20 sulfate per dosage unit.

21 (c) *Stimulants.* —

22 Pyrovalerone.

23 (d) Any compound, mixture, or preparation containing as its single active ingredient
24 ephedrine, pseudoephedrine, or phenylpropanolamine, their salts or optical isomers, or salts of
25 optical isomers except products which are for pediatric use primarily intended for administration
26 to children under the age of 12: *Provided*, That neither the offenses set forth in §60A-4-401 nor
27 the penalties therein shall be applicable to ephedrine, pseudoephedrine, or phenylpropanolamine
28 which shall be subject to the provisions of §60A-10-1 *et seq.* of this code.

29 (e) *Depressants.* —

30 Ezogabine [N-[2-amino-4-(9-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester];

31 Lacosamide [(R)-2-acetoamido- N -benzyl-3-methoxy-propionamide];

32 ~~Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid]; and~~

33 Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl] butanamide) (also referred to as
34 BRV; UCB-34714; Briviact).

35 (f) *Other substances:* —

36 Gabapentin

37 Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid];

38 Epidiolex (contains cannabidiol (CBD)).

