

WEST VIRGINIA LEGISLATURE

2020 REGULAR SESSION

Engrossed

Senate Bill 847

BY SENATORS AZINGER, BALDWIN, CLEMENTS, CLINE,
HARDESTY, JEFFRIES, LINDSAY, MAYNARD,
PITSENBARGER, ROMANO, RUCKER, SMITH, WELD,
WOELFEL, AND TRUMP

[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- 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 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- piperidyl]-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 Hydromorphenol;
- 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 3,4,5-trimethoxy amphetamine;

- 124 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
- 125 Alpha-methyltryptamine (other name: AMT);
- 126 Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole;
- 127 3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N- dimethyltryptamine;
- 128 mappine;
- 129 Diethyltryptamine; some trade and other names: N, N-Diethyltryptamine; DET;
- 130 Dimethyltryptamine; some trade or other names: DMT;
- 131 5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);
- 132 Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-
- 133 methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;
- 134 Lysergic acid diethylamide;
- 135 ~~Marihuana~~ Marijuana (Cannabis, sp.);
- 136 Mescaline;
- 137 Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,
- 138 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;
- 139 Peyote; meaning all parts of the plant presently classified botanically as *Lophophora*
- 140 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from any part of such
- 141 plant, and every compound, manufacture, salts, immediate derivative, mixture, or preparation of
- 142 such plant, its seeds, or extracts;
- 143 N-ethyl-3-piperidyl benzilate;
- 144 N-methyl-3-piperidyl benzilate;
- 145 Psilocybin;
- 146 Psilocyn;
- 147 Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or
- 148 in the resinous extractives of *Cannabis, sp.* and/or synthetic substances, immediate derivatives

149 and their isomers with similar chemical structure and pharmacological activity such as the
150 following:

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

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

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

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

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

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

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

163 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy;

164 4-methylmethcathinone (Mephedrone);

165 3,4-methylenedioxypropylvalerone (MDPV);

166 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

167 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

168 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

169 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

170 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

171 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

172 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

173 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);

174 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);

175 3,4-Methylenedioxy-N-methylcathinone (Methylone);
176 2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, its optical isomers, salts and
177 salts of isomers;
178 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-
179 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);
180 Alpha-methyltryptamine (other name: AMT);
181 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
182 Synthetic Cannabinoids as follows:
183 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol {also known as CP
184 47,497 and homologues};
185 rel-2-[(1S,3R)-3-hydroxycyclohexyl]-5-(2-methylnonan-2-yl)phenol {also known as CP
186 47,497-C8 homolog};
187 [(6aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
188 tetrahydrobenzo[c]chromen-1-ol] {also known as HU-210};
189 (dexanabinol);
190 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
191 tetrahydrobenzo[c]chromen-1-ol {also known as HU-211};
192 1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};
193 1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};
194 (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-015};
195 (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};
196 [1-[2-(4-morpholinyl)ethyl]-1H-indol-3-yl]-1-naphthalenyl-methanone {also known as
197 JWH-200};
198 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};
199 2-((1S,2S,5S)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl)-5-(2-methyloctan-2-yl)phenol
200 {also known as CP 55,940};

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

227 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
228 (ADB-FUBINACA);

229 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
230 (MDMB-CHMICA);

231 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate
232 (MDMB-FUBINACA);

233 Tetrahydrocannabinols:

234 DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.

235 DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.

236 DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.

237 Synthetic Phenethylamines:

238 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
239 NBOMe);

240 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-
241 NBOMe);

242 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
243 NBOMe);

244 Synthetic Opioids (~~including~~ including their isomers, esters, ethers, salts, and salts of
245 isomers, esters, and ethers):

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

247 furanyl fentanyl;

248 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
249 47700);

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

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

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

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

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

258 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide

259 (also known as U-48800);

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

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

264 *Opioid Receptor Agonist.* —

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

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

270 JWH 015;

271 JWH 018;

272 JWH 019;

273 JWH 073;

274 JWH 081;

275 JWH 122;

276 JWH 200;

277 JWH 210;

278 JWH 398;

279 AM 2201;

280 WIN 55,212.

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

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

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

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

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

298 JWH 250;

299 JWH 203;

300 JWH 251;

301 JWH 302.

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

305 CP 47,497 and its homologues and analogs;

306 Cannabicyclohexanol;

307 CP 55,940.

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

312 AM 694;

313 Pravadoline WIN 48,098;

314 RCS 4;

315 AM 679.

316 [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-
317 naphthalenemethanone. This shall include WIN 55,212-2.

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

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

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

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

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

335 Tryptamines:

336 5-methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT);

337 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT);

338 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT);

339 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);

340 4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT);

341 5-methoxy- α -methyltryptamine (5-MeO-AMT);

342 4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT);

343 4-hydroxy Diethyltryptamine (4-HO-DET);

344 5-methoxy- N,N- diallyltryptamine (5-MeO-DALT);

345 4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT);

346 ~~4-hydroxy Diethyltryptamine (4-HO-DET)~~

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

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

349 5-Fluoro-MN-24 (1-(5-Fluoropentyl)-N-(naphthalen-1-yl)-1H-indole-3-carboxamide);

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

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

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

353 Methyl-Ethylaminopentiophenone;

- 354 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
- 355 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
- 356 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
- 357 carboxamide);
- 358 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
- 359 methylbutanoat);
- 360 MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
- 361 SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
- 362 SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
- 363 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
- 364 Methyl-Ethylaminopentiophenone;
- 365 FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
- 366 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
- 367 5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-
- 368 3- carboxamide);
- 369 MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
- 370 methylbutanoat);
- 371 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
- 372 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-
- 373 a][1,4]benzodiazepine);
- 374 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
- 375 benzodiazepin-2-one);
- 376 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
- 377 a][1,4]diazepine);
- 378 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
- 379 a][1,4]benzodiazepine);

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

381 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-

382 a][1,4]benzodiazepine);

383 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzof[1,2,4]triazolo[4,3-

384 a][1,4]diazepine);

385 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-

386 one) ;

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

388 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-

389 a][1,4]benzodiazepine).

390 (e) *Depressants*. —

391 4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-

392 carboxamide);

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

394 2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-cyclohexanone,

395 monohydrochloride);

396 4-MEAP (2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one);

397 Mecloqualone;

398 Methaqualone;

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

400 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3

401 a][1,4]benzodiazepine);

402 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-

403 benzodiazepin-2-one);

404 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-

405 a][1,4]diazepine);

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

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

458 3-Fluorophenmetrazine (2-(3-Fluorophenyl)-3-methylmorpholine);

459 (g) Temporary listing of substances subject to emergency scheduling. Any material,
460 compound, mixture, or preparation which contains any quantity of the following substances:

461 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts,
462 and salts of isomers;

463 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
464 isomers, salts and salts of isomers;

465 N-benzylpiperazine, also known as BZP;

466 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);

467 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-
468 butyramide);

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

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

472 3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-
473 phenylbutyramide);

474 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
475 yl)butyramide);

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

477 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
478 carboxamide);

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

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

481 Synthetic cathinones or any compound, except bupropion or compounds listed under a
482 different schedule, or compounds used within legitimate and approved medical research,
483 structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic

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

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

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

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

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

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