

WEST VIRGINIA LEGISLATURE

2017 REGULAR SESSION

ENROLLED

Committee Substitute

for

House Bill 2526

BY DELEGATES ELLINGTON, SUMMERS, SOBONYA AND

ROHRBACH

[Passed April 8, 2017; in effect ninety days from passage.]

1 AN ACT to amend and reenact §60A-2-201, §60A-2-204, §60A-2-206, §60A-2-210 and §60A-2-
2 212 of the Code of West Virginia, 1931, as amended, all relating to classifying additional
3 drugs to Schedules I, II, IV and V of controlled substances; and adding a provision relating
4 to the scheduling of a cannabidiol in a product approved by the Food and Drug
5 Administration.

Be it enacted by the Legislature of West Virginia:

1 That §60A-2-201, §60A-2-204, §60A-2-206, §60A-2-210 and §60A-2-212 of the Code of
2 West Virginia, 1931, as amended, be amended and reenacted, all to read as follows:

ARTICLE 2. STANDARDS AND SCHEDULES.

§60A-2-201. Authority of state Board of Pharmacy; recommendations to Legislature.

1 (a) The state Board of Pharmacy shall administer the provisions of this chapter. It shall
2 also, on the first day of each regular legislative session, recommend to the Legislature which
3 substances should be added to or deleted from the schedules of controlled substances contained
4 in this article or reschedule therein. The state Board of Pharmacy shall also have the authority
5 between regular legislative sessions, on an emergency basis, to add to or delete from the
6 schedules of controlled substances contained in this article or reschedule such substances based
7 upon the recommendations and approval of the federal food, drug and cosmetic agency, and shall
8 report such actions on the first day of the regular legislative session immediately following said
9 actions.

10 In making any such recommendation regarding a substance, the state Board of Pharmacy
11 shall consider the following factors:

- 12 (1) The actual or relative potential for abuse;
- 13 (2) The scientific evidence of its pharmacological effect, if known;
- 14 (3) The state of current scientific knowledge regarding the substance;
- 15 (4) The history and current pattern of abuse;
- 16 (5) The scope, duration and significance of abuse;

17 (6) The potential of the substance to produce psychic or physiological dependence liability;
18 and

19 (7) Whether the substance is an immediate precursor of a substance already controlled
20 under this article.

21 (b) After considering the factors enumerated in subsection (a), the state Board of
22 Pharmacy shall make findings with respect to the substance under consideration. If it finds that
23 any substance not already controlled under any schedule has a potential for abuse, it shall
24 recommend to the Legislature that the substance be added to the appropriate schedule. If it finds
25 that any substance already controlled under any schedule should be rescheduled or deleted, it
26 shall so recommend to the Legislature.

27 (c) If the state Board of Pharmacy designates a substance as an immediate precursor,
28 substances which are precursors of the controlled precursor shall not be subject to control solely
29 because they are precursors of the controlled precursor.

30 (d) If any substance is designated, rescheduled or deleted as a controlled substance under
31 federal laws and notice thereof is given to the state Board of Pharmacy, the board shall
32 recommend similar control of such substance to the Legislature, specifically stating that such
33 recommendation is based on federal action and the reasons why the federal government deemed
34 such action necessary and proper.

35 (e) The authority vested in the board by subsection (a) of this section shall not extend to
36 distilled spirits, wine, malt beverages or tobacco as those terms are defined or used in other
37 chapters of this code nor to any nonnarcotic substance if such substance may under the "Federal
38 Food, Drug and Cosmetic Act" and the law of this state lawfully be sold over the counter without
39 a prescription.

40 (f) Notwithstanding any provision of this chapter to the contrary, the sale, wholesale,
41 distribution or prescribing of a cannabidiol in a product approved by the Food and Drug
42 Administration is permitted and shall be placed on the schedule as provided for by the Drug
43 Enforcement Administration.

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.

3 (b) Opiates. Unless specifically excepted or unless listed in another schedule, any of the
4 following opiates, including their isomers, esters, ethers, salts and salts of isomers, esters and
5 ethers, whenever the existence of such isomers, esters, ethers and salts is possible within the
6 specific chemical designation (for purposes of subdivision (34) of this subsection only, the term
7 isomer includes the optical and geometric isomers):

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

10 (2) Acetylmethadol;

11 (3) Allylprodine;

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

14 (5) Alphameprodine;

15 (6) Alphamethadol;

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

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

20 (9) Benzethidine;

21 (10) Betacetylmethadol;

22 (11) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl) -4- piperidiny]-N-
23 phenylpropanamide);

24 (12) Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2- hydroxy-2-phenethyl)-3-methyl-
25 4-piperidiny]-N-phenylpropanamide);

- 26 (13) Betameprodine;
- 27 (14) Betamethadol;
- 28 (15) Betaprodine;
- 29 (16) Clonitazene;
- 30 (17) Dextromoramide;
- 31 (18) Diampromide;
- 32 (19) Diethylthiambutene;
- 33 (20) Difenoxin;
- 34 (21) Dimenoxadol;
- 35 (22) Dimepheptanol;
- 36 (23) Dimethylthiambutene;
- 37 (24) Dioxaphetyl butyrate;
- 38 (25) Dipipanone;
- 39 (26) Ethylmethylthiambutene;
- 40 (27) Etonitazene;
- 41 (28) Etoxidine;
- 42 (29) Furethidine;
- 43 (30) Hydroxypethidine;
- 44 (31) Ketobemidone;
- 45 (32) Levomoramide;
- 46 (33) Levophenacymorphan;
- 47 (34) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4- piperidyl]-N-
- 48 phenylpropanamide);
- 49 (35) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4- piperidiny]—
- 50 phenylpropanamide);
- 51 (36) Morpheridine;

- 52 (37) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- 53 (38) Noracymethadol;
- 54 (39) Norlevorphanol;
- 55 (40) Normethadone;
- 56 (41) Norpipanone;
- 57 (42) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl]
- 58 propanamide);
- 59 (43) PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 60 (44) Phenadoxone;
- 61 (45) Phenampromide;
- 62 (46) Phenomorphan;
- 63 (47) Phenoperidine;
- 64 (48) Piritramide;
- 65 (49) Proheptazine;
- 66 (50) Properidine;
- 67 (51) Propiram;
- 68 (52) Racemoramide;
- 69 (53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl) ethyl-4- piperidinyl]-propanamide);
- 70 (54) Tilidine;
- 71 (55) Trimeperidine.

72 (c) *Opium derivatives*. — Unless specifically excepted or unless listed in another
73 schedule, any of the following opium immediate derivatives, its salts, isomers and salts of isomers
74 whenever the existence of such salts, isomers and salts of isomers is possible within the specific
75 chemical designation:

- 76 (1) Acetorphine;
- 77 (2) Acetyldihydrocodeine;

- 78 (3) Benzylmorphine;
- 79 (4) Codeine methylbromide;
- 80 (5) Codeine-N-Oxide;
- 81 (6) Cyprenorphine;
- 82 (7) Desomorphine;
- 83 (8) Dihydromorphine;
- 84 (9) Drotebanol;
- 85 (10) Etorphine (except HCl Salt);
- 86 (11) Heroin;
- 87 (12) Hydromorphenol;
- 88 (13) Methyldesorphine;
- 89 (14) Methyldihydromorphine;
- 90 (15) Morphine methylbromide;
- 91 (16) Morphine methylsulfonate;
- 92 (17) Morphine-N-Oxide;
- 93 (18) Myrophine;
- 94 (19) Nicocodeine;
- 95 (20) Nicomorphine;
- 96 (21) Normorphine;
- 97 (22) Pholcodine;
- 98 (23) Thebacon.

99 (d) *Hallucinogenic substances*. — Unless specifically excepted or unless listed in another
100 schedule, any material, compound, mixture or preparation, which contains any quantity of the
101 following hallucinogenic substances, or which contains any of its salts, isomers and salts of
102 isomers, whenever the existence of such salts, isomers, and salts of isomers is possible within

103 the specific chemical designation (for purposes of this subsection only, the term “isomer” includes
104 the optical, position and geometric isomers):

105 (1) Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-
106 1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; alpha-ET; and AET;

107 (2) 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
108 dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;

109 (3) 4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-
110 dimethoxyphenyl)-1-aminoethane; alpha- desmethyl DOB; 2C-B, Nexus;

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

113 (B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe).

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

115 (5) 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-
116 methylphenethylamine; 2,5-DMA;

117 (6) 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

118 (7) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

119 (8) 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-
120 methylphenethylamine; paramethoxyamphetamine; PMA;

121 (9) 5-methoxy-3, 4-methylenedioxy-amphetamine;

122 (10) 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-
123 dimethoxy-alpha-methylphenethylamine; “DOM”; and “STP”;

124 (11) 3,4-methylenedioxy amphetamine;

125 (12) 3,4-methylenedioxymethamphetamine (MDMA);

126 (13) 3,4-methylenedioxy-N-ethylamphetamine (also known as – ethyl-alpha-methyl-3,4
127 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

- 128 (14) N-hydroxy-3,4-methylenedioxyamphetamine (also known as – hydroxy-alpha-methyl-
129 3,4 (methylenedioxy) phenethylamine, and – hydroxy MDA);
- 130 (15) 3,4,5-trimethoxy amphetamine;
- 131 (16) 5-methoxy-N, N-dimethyltryptamine (5-MeO-DMT);
- 132 (17) Alpha-methyltryptamine (other name: AMT);
- 133 (18) Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-
134 hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-
135 dimethyltryptamine; mappine;
- 136 (19) Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;
- 137 (20) Dimethyltryptamine; some trade or other names: DMT;
- 138 (21) 5-Methoxy-N, N-diisopropyltryptamine (5-MeO-DIPT);
- 139 (22) Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-
140 octahydro-2-methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe
141 iboga;
- 142 (23) Lysergic acid diethylamide;
- 143 (24) Marijuana;
- 144 (25) Mescaline;
- 145 (26) Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-
146 tetrahydro-6, 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;
- 147 (27) Peyote; meaning all parts of the plant presently classified botanically as *Lophophora*
148 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from any part of such
149 plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of
150 such plant, its seeds or extracts;
- 151 (28) N-ethyl-3-piperidyl benzilate;
- 152 (29) N-methyl-3-piperidyl benzilate;
- 153 (30) Psilocybin;

154 (31) Psilocyn;

155 (32) Tetrahydrocannabinols; synthetic equivalents of the substances contained in the
156 plant, or in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate
157 derivatives and their isomers with similar chemical structure and pharmacological activity such as
158 the following:

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

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

161 delta-3,4 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 (33) Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-
165 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,
166 cyclohexamine, PCE;

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

169 (35) Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-
170 cyclohexyl]-piperidine, 2-thienyl analog of phencyclidine; TPCP, TCP;

171 (36) 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy.

172 (37) 4-methylmethcathinone (Mephedrone);

173 (38) 3,4-methylenedioxypropylvalerone (MDPV);

174 (39) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

175 (40) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

176 (41) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

177 (42) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

178 (43) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

179 (44) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

- 180 (45) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);
- 181 (46) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N);
- 182 (47) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);
- 183 (48) 3,4-Methylenedioxy-N-methylcathinone (Methylone);
- 184 (49) (2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, itsoptical isomers, salts
- 185 and salts of isomers
- 186 (50) 5-methoxy-N, N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-
- 187 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);
- 188 (51) Alpha-methyltryptamine (other name: AMT);
- 189 (52) 5-methoxy-N, N-diisopropyltryptamine (other name: 5-MeO-DIPT);
- 190 (53) Synthetic Cannabinoids as follows:
- 191 (A) 2-[(1R,3S)-3-hydroxycyclohexyl]-5- (2-methyloctan-2-yl) phenol {also known as CP
- 192 47,497 and homologues};
- 193 (B) rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl) phenol {also known as
- 194 CP 47,497-C8 homolog};
- 195 (C) [(6aR)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7,10,10a-
- 196 tetrahydrobenzo[c]chromen-1-ol]] {also known as HU-210};
- 197 (D) (dexanabinol);
- 198 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
- 199 tetrahydrobenzol[c]chromen-1-ol) {also known as HU-211};
- 200 (E) 1-Pentyl-3-(1-naphthoyl) indole {also known as JWH-018};
- 201 (F) 1-Butyl-3-(1-naphthoyl) indole {also known as JWH-073};
- 202 (G) (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-
- 203 015};
- 204 (H) (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};

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

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

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

210 (L) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-
211 122};

212 (M) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-
213 398};

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

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

217 (P) 1-pentyl-3-[1-(4-methoxynaphthoyl) indole (JWH-081);

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

219 (R) 1-(5-fluoropentyl)-3-(2-iodobenzoyl) indole (AM694).

220 (54) Synthetic cannabinoids or any material, compound, mixture or preparation which
221 contains any quantity of the following substances, including their analogues, congeners,
222 homologues, isomers, salts and salts of analogues, congeners, homologues and isomers, as
223 follows:

224 (A) CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
225 YL) phenol);

226 (B) HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-
227 6A,7,10, 10A-tetrahydrobenzo[C] chromen-1-OL)];

228 (C) HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-
229 methyloctan-2-YL)-6A,7,10,10 atetrahydrobenzo [C] chromen-1-OL);

230 (D) JWH-018, 1-pentyl-3-(1-naphthoyl) indole;

231 (E) JWH-019, 1-hexyl-3-(1-naphthoyl) indole;

232 (F) JWH-073, 1-butyl-3-(1-naphthoyl) indole;

233 (G) JWH-200, (1-(2-morpholin-4-ylethyl) indol-3-yl)- Naphthalen-1-ylmethanone;

234 (H) JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl) indole.

235 (55) Synthetic cannabinoids including any material, compound, mixture or preparation that
236 is not listed as a controlled substance in Schedule I through V, is not a federal Food and Drug
237 Administration approved drug or used within legitimate and approved medical research and which
238 contains any quantity of the following substances, their salts, isomers, whether optical positional
239 or geometric, analogues, homologues and salts of isomers, analogues and homologues, unless
240 specifically exempted, whenever the existence of these salts, isomers, analogues, homologues
241 and salts of isomers, analogues and homologues if possible within the specific chemical
242 designation:

243 (A) Tetrahydrocannabinols meaning tetrahydrocannabinols which are naturally contained
244 in a plant of the genus cannabis as well as synthetic equivalents of the substances contained in
245 the plant or in the resinous extractives of cannabis or synthetic substances, derivatives and their
246 isomers with analogous chemical structure and or pharmacological activity such as the following:

247 (i) DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.

248 (ii) DELTA-6 CIS OR trans tetrahydrocannabinol and their Optical isomers.

249 (iii) DELTA-3,4 CIS OR their trans tetrahydrocannabinol and their optical isomers.

250 (B) Naphthoyl indoles or any compound containing a 3-(-1- Naphthoyl) indole structure with
251 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole
252 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall
253 include the following:

254 (i) JWH 015;

255 (ii) JWH 018;

256 (iii) JWH 019;

257 (iv) JWH 073;

258 (v) JWH 081;

259 (vi) JWH 122;

260 (vii) JWH 200;

261 (viii) JWH 210;

262 (ix) JWH 398;

263 (x) AM 2201;

264 (xi) WIN 55,212.

265 (56) Synthetic Phenethylamines (including their optical, positional, and geometric isomers,
266 salts and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers):

267 (A) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
268 NBOMe);

269 (B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-
270 C-NBOMe);

271 (C) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/
272 2C-B-NBOMe);

273 (57) Synthetic Opioids (including their isomers, esters, ethers, salts and salts of isomers,
274 esters and ethers):

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

276 (B) furanyl fentanyl;

277 (C) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
278 47700);

279 (D) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
280 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);

281 (E) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also
282 known as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-
283 hydroxythiofentanyl).

284 (58) Opioid Receptor Agonist (including its isomers, esters, ethers, salts, and salts of
285 isomers, esters and ethers):

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

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

291 (60) Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure
292 with substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the
293 pyrrole 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 147 and JWH 307.

295 (61) Naphthylmethylindenenes or any compound containing a Naphthylideneindene
296 structure with substitution at the 3- Position of the indene ring whether or not further substituted
297 in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent.
298 This shall include, but not be limited to, JWH 176.

299 (62) Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure
300 with substitution at the nitrogen atom of the indole ring whether or not further substituted in the
301 indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall
302 include the following:

303 (A) RCS-8, SR-18 OR BTM-8;

304 (B) JWH 250;

305 (C) JWH 203;

306 (D) JWH 251;

307 (E) JWH 302.

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

311 (A) CP 47,497 and its homologues and analogs;

312 (B) Cannabicyclohexanol;

313 (C) CP 55,940.

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

318 (A) AM 694;

319 (B) Pravadoline WIN 48,098;

320 (C) RCS 4;

321 (D) AM 679.

322 (65) [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-
323 YL]-1-napthalenymethanone. This shall include WIN 55,212-2.

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

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

330 (68) Tetramethylcyclopropylindoles or any compound containing A 3-
331 tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring

332 whether or not further substituted in the indole ring to any extent and whether or not substituted
333 in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.

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

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

341 (71) Tryptamines:

342 (A) 5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)

343 (B) 4-hydroxy-N, N-diisopropyltryptamine (4-HO-DiPT)

344 (C) 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)

345 (D) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)

346 (E) 4-acetoxy-N, N-diisopropyltryptamine (4-AcO-DiPT)

347 (F) 5-methoxy- α -methyltryptamine (5-MeO-AMT)

348 (G) 4-methoxy-N, N-Dimethyltryptamine (4-MeO-DMT)

349 (H) 4-hydroxy Diethyltryptamine (4-HO-DET)

350 (I) 5- methoxy- N, N- diallyltryptamine (5-MeO-DALT)

351 (J) 4-acetoxy-N, N-Dimethyltryptamine (4-AcO DMT)

352 (K) 4-hydroxy Diethyltryptamine (4-HO-DET)

353 (72) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-
354 carboxamide (AB-CHMINACA);

355 (73) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (AB-
356 PINACA);

357 (74) [1-(5-fluoropentyl)-1H-indazol-3-yl (naphthalen-1-yl)methanone (THJ-2201);

358 (75) quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB-22; QUPIC);

359 (76) quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22);

360 (77) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-

361 carboxamide (AB-FUBINACA);

362 (78) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide

363 (ADB-PINACA); and

364 (79) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-

365 carboxamide (common names, MAB-CHMINACA and ADB-CHMINACA);

366 (e) *Depressants*. — Unless specifically excepted or unless listed in another schedule, any
367 material, compound, mixture, or preparation which contains any quantity of the following
368 substances having a depressant effect on the central nervous system, including its salts, isomers
369 and salts of isomers whenever the existence of such salts, isomers and salts of isomers is
370 possible within the specific chemical designation:

371 (1) Mecloqualone;

372 (2) Methaqualone.

373 (f) *Stimulants*. — Unless specifically excepted or unless listed in another schedule, any
374 material, compound, mixture, or preparation which contains any quantity of the following
375 substances having a stimulant effect on the central nervous system, including its salts, isomers
376 and salts of isomers:

377 (1) Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
378 dihydro-5-phenyl-2-oxazolamine;

379 (2) Cathinone; some trade or other names: 2-amino-1-phenyl-1- propanone, alpha-
380 aminopropiophenone, 2-aminopropiophenone and norephedrone;

381 (3) Fenethylamine;

382 (4) Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
383 isomers and salts of optical isomers; some other names: (2-(methylamino)-propiofenone; alpha-

384 (methylamino)propiofenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—
385 methylaminopropiofenone; monomethylpropion; 3,4-methylenedioxypropiofenone and/or
386 mephedrone; 3,4-methylenedioxypropiofenone (MPVD); ephedrone; N-methylcathinone;
387 methylcathinone; AL-464; AL-422; AL- 463 and UR1432;

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

389 (6) N-ethylamphetamine;

390 (7) N,N-dimethylamphetemine; also known as N,N-alpha- trimethyl-benzeneethanamine;
391 N,N-alpha-trimethylphenethylamine.

392 (8) Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and
393 salts of isomers.

394 (9) Substituted amphetamines:

395 (A) 2-Fluoroamphetamine

396 (B) 3-Fluoroamphetamine

397 (C) 4-Fluoroamphetamine

398 (D) 2-chloroamphetamine

399 (E) 3-chloroamphetamine

400 (F) 4-chloroamphetamine

401 (G) 2-Fluoromethamphetamine

402 (H) 3-Fluoromethamphetamine

403 (I) 4-Fluoromethamphetamine

404 (J) 4-chloromethamphetamine

405 (10) 4-methyl-N-ethylcathinone (4-MEC);

406 (11) 4-methyl-alpha-pyrrolidinopropiofenone (4-MePPP);

407 (12) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone);

408 (13) 2-(methylamino)-1-phenylpentan-1-one (pentedrone);

409 (14) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone);

410 (15) 4-fluoro-N-methylcathinone (4-FMC);

411 (16) 3-fluoro-N-methylcathinone (3-FMC);

412 (17) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); and

413 (18) Alpha-pyrrolidinobutiophenone (α -PBP).

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

416 (1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers,
417 salts, and salts of isomers.

418 (2) N-[1-(2-thienyl) methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
419 isomers, salts and salts of isomers.

420 (3) N-benzylpiperazine, also known as BZP.

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

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

427 (A) By substitution in the ring system to any extent with Alkyl, alkylendioxy, alkoxy,
428 haloalkyl, hydroxyl or halide Substituents whether or not further substituted in the ring system by
429 one or more other univalent substituents.

430 (B) By substitution at the 3-Position with an acyclic alkyl substituent.

431 (C) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl or
432 methoxybenzyl groups.

433 (D) By inclusion of the 2-amino nitrogen atom in a cyclic structure.

434 (2) Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist
435 as demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV

436 and V, not federal Food and Drug Administration approved drug or used within legitimate,
437 approved medical research.

§60A-2-206. Schedule II.

1 (a) Schedule II consists 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.

3 (b) *Substances, vegetable origin or chemical synthesis.* — Unless specifically excepted or
4 unless listed in another schedule, any of the following substances whether produced directly or
5 indirectly by extraction from substances of vegetable origin, or independently by means of
6 chemical synthesis, or by a combination of extraction and chemical synthesis:

7 (1) Opium and opiate, and any salt, compound, derivative or preparation of opium or opiate
8 excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmeferene,
9 naloxone and naltrexone, and their respective salts, but including the following:

10 (A) Raw opium;

11 (B) Opium extracts;

12 (C) Opium fluid;

13 (D) Powdered opium;

14 (E) Granulated opium;

15 (F) Tincture of opium;

16 (G) Codeine;

17 (H) Dihydroetorphine;

18 (I) Ethylmorphine;

19 (J) Etorphine hydrochloride;

20 (K) Hydrocodone;

21 (L) Hydromorphone;

22 (M) Metopon;

23 (N) Morphine;

24 (O) Oripavine;

25 (P) Oxycodone;

26 (Q) Oxymorphone; and

27 (R) Thebaine;

28 (2) Any salt, compound, derivative or preparation thereof which is chemically equivalent
29 or identical with any of the substances referred to in subdivision (1) of this subsection, except that
30 these substances shall not include the isoquinoline alkaloids of opium;

31 (3) Opium poppy and poppy straw;

32 (4) Coca leaves and any salt, compound, derivative or preparation of coca leaves
33 (including cocaine and ecgonine and their salts, isomers, derivatives and salts of isomers and
34 derivatives), and any salt, compound, derivative or preparation thereof which is chemically
35 equivalent or identical with any of these substances, except that the substances shall not include
36 decocainized coca leaves or extractions of coca leaves, which extractions do not contain cocaine
37 or ecgonine;

38 (5) Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid or
39 powder form which contains the phenanthrene alkaloids of the opium poppy).

40 (c) *Opiates*. — Unless specifically excepted or unless in another schedule, any of the
41 following opiates, including its isomers, esters, ethers, salts and salts of isomers, esters and
42 ethers whenever the existence of such isomers, esters, ethers and salts is possible within the
43 specific chemical designation, dextrorphan and levopropoxyphene excepted:

44 (1) Alfentanil;

45 (2) Alphaprodine;

46 (3) Anileridine;

47 (4) Bezitramide;

48 (5) Bulk dextropropoxyphene (nondosage forms);

49 (6) Carfentanil;

- 50 (7) Dihydrocodeine;
- 51 (8) Diphenoxylate;
- 52 (9) Fentanyl;
- 53 (10) Isomethadone;
- 54 (11) Levo-alphaacetylmethadol; some other names: levo-alpha-acetylmethadol,
- 55 levomethadyl acetate, LAAM;
- 56 (12) Levomethorphan;
- 57 (13) Levorphanol;
- 58 (14) Metazocine;
- 59 (15) Methadone;
- 60 (16) Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane;
- 61 (17) Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic acid;
- 62 (18) Pethidine; (meperidine);
- 63 (19) Pethidine-Intermediate-A, 4-cyano-1-methyl-4- phenylpiperidine;
- 64 (20) Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;
- 65 (21) Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;
- 66 (22) Phenazocine;
- 67 (23) Piminodine;
- 68 (24) Racemethorphan;
- 69 (25) Racemorphan;
- 70 (26) Remifentanil;
- 71 (27) Sufentanil;
- 72 (28) Tapentadol;

73 (29) Thiafentanil (4-(methoxycarbonyl)-4-(N-phenmethoxyacetamido)-1-2-
74 (thienyl)ethylpiperidine), including its isomers, esters, ethers, salts and salts of isomers, esters
75 and ethers.

76 (d) *Stimulants*. — Unless specifically excepted or unless listed in another schedule, any
77 material, compound, mixture or preparation which contains any quantity of the following
78 substances having a stimulant effect on the central nervous system:

79 (1) Amphetamine, its salts, optical isomers and salts of its optical isomers;

80 (2) Methamphetamine, its salts, isomers and salts of its isomers;

81 (3) Methylphenidate;

82 (4) Phenmetrazine and its salts; and

83 (5) Lisdexamfetamine.

84 (e) *Depressants*. — Unless specifically excepted or unless listed in another schedule, any
85 material, compound, mixture or preparation which contains any quantity of the following
86 substances having a depressant effect on the central nervous system, including its salts, isomers
87 and salts of isomers whenever the existence of such salts, isomers and salts of isomers is
88 possible within the specific chemical designation:

89 (1) Amobarbital;

90 (2) Glutethimide;

91 (3) Pentobarbital;

92 (4) Phencyclidine;

93 (5) Secobarbital.

94 (f) *Hallucinogenic substances*:

95 Nabilone: [Another name for nabilone: (+-)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8, 10,
96 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9H-dibenzo [b,d] pyran-9-one].

97 (g) *Immediate precursors*. — Unless specifically excepted or unless listed in another
98 schedule, any material, compound, mixture, or preparation which contains any quantity of the
99 following substances:

- 100 (1) Immediate precursor to amphetamine and methamphetamine:
101 (A) Phenylacetone;
102 (B) Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl
103 benzyl ketone;
- 104 (2) Immediate precursors to phencyclidine (PCP):
105 (A) 1-phenylcyclohexylamine; and
106 (B) 1-piperidinocyclohexanecarbonitrile (PCC).
- 107 (3) Immediate precursor to fentanyl:
108 4-anilino-N-phenethyl-4-piperidine (ANPP).

§60A-2-210. Schedule IV.

- 1 (a) Schedule IV 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.
- 3 (b) *Narcotic drugs.* — Unless specifically excepted or unless listed in another schedule,
4 any material, compound, mixture or preparation containing any of the following narcotic drugs, or
5 their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth
6 below:
- 7 (1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine
8 sulfate per dosage unit;
- 9 (2) Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-
10 propionoxybutane).
- 11 (c) *Depressants.* — Unless specifically excepted or unless listed in another schedule, any
12 material, compound, mixture or preparation which contains any quantity of the following
13 substances, including its salts, isomers and salts of isomers whenever the existence of such salts,
14 isomers and salts of isomers is possible within the specific chemical designation:
- 15 (1) Alprazolam;
16 (2) Barbital;

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- 17 (3) Bromazepam;
- 18 (4) Camazepam;
- 19 (5) Carisoprodol;
- 20 (6) Chloral betaine;
- 21 (7) Chloral hydrate;
- 22 (8) Chlordiazepoxide;
- 23 (9) Clobazam;
- 24 (10) Clonazepam;
- 25 (11) Clorazepate;
- 26 (12) Clotiazepam;
- 27 (13) Cloxazolam;
- 28 (14) Delorazepam;
- 29 (15) Diazepam;
- 30 (16) Dichloralphenazone;
- 31 (17) Estazolam;
- 32 (18) Ethchlorvynol;
- 33 (19) Ethinamate;
- 34 (20) Ethyl loflazepate;
- 35 (21) Fludiazepam;
- 36 (22) Flunitrazepam;
- 37 (23) Flurazepam;
- 38 (24) Fospropofol;
- 39 (25) Halazepam;
- 40 (26) Haloxazolam;
- 41 (27) Ketazolam;
- 42 (28) Loprazolam;

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- 43 (29) Lorazepam;
- 44 (30) Lormetazepam;
- 45 (31) Mebutamate;
- 46 (32) Medazepam;
- 47 (33) Meprobamate;
- 48 (34) Methohexital;
- 49 (35) Methylphenobarbital (mephobarbital);
- 50 (36) Midazolam;
- 51 (37) Nimetazepam;
- 52 (38) Nitrazepam;
- 53 (39) Nordiazepam;
- 54 (40) Oxazepam;
- 55 (41) Oxazolam;
- 56 (42) Paraldehyde;
- 57 (43) Petrichloral;
- 58 (44) Phenobarbital;
- 59 (45) Pinazepam;
- 60 (46) Prazepam;
- 61 (47) Quazepam;
- 62 (48) Temazepam;
- 63 (49) Tetrazepam;
- 64 (50) Triazolam;
- 65 (51) Zaleplon;
- 66 (52) Zolpidem;
- 67 (53) Zopiclone'

68 (54) Suvorexant ([[(7R)-4-(5-chloro-1,3-benzoxazol-2-yl)-7-methyl-1,4-diazepan-1-yl] [5-
69 methyl-2-(2H-1,2,3-triazol-2-yl)phenyl]methanone).

70 (d) Any material, compound, mixture or preparation which contains any quantity of the
71 following substance, including its salts, isomers (whether optical, position or geometric) and salts
72 of such isomers whenever the existence of such salts, isomers and salts of isomers is possible:
73 Fenfluramine and Dexfenfluramine.

74 (e) *Stimulants*. — Unless specifically excepted or unless listed in another schedule, any
75 material, compound, mixture or preparation which contains any quantity of the following
76 substances having a stimulant effect on the central nervous system, including its salts, isomers
77 and salts of isomers:

- 78 (1) Cathine ((+)-norpseudoephedrine);
- 79 (2) Diethylpropion;
- 80 (3) Fencamfamin;
- 81 (4) Fenproporex;
- 82 (5) Mazindol;
- 83 (6) Mefenorex;
- 84 (7) Modafinil;
- 85 (8) Pemoline (including organometallic complexes and chelates thereof);
- 86 (9) Phentermine;
- 87 (10) Pipradrol;
- 88 (11) Sibutramine;
- 89 (12) SPA ((-)-1-dimethylamino-1,2-diphenylethane);
- 90 (13) Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropyl
91 [(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid);

92 (f) *Other substances*. — Unless specifically excepted or unless listed in another schedule,
93 any material, compound, mixture or preparation which contains any quantity of the following
94 substances, including its salts:

95 (1) Pentazocine;

96 (2) Butorphanol;

97 (3) Tramadol (2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol).

98 Amyl nitrite, butyl nitrite, isobutyl nitrite and the other organic nitrites are controlled
99 substances and no product containing these compounds as a significant component shall be
100 possessed, bought or sold other than pursuant to a bona fide prescription or for industrial or
101 manufacturing purposes.

§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.

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

9 (1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;

10 (2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams;

11 (3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams;

12 (4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of
13 atropine sulfate per dosage unit;

14 (5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams;

15 (6) Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine
16 sulfate per dosage unit.

17 (c) *Stimulants*. — Unless specifically exempted or excluded or unless listed in another
18 schedule, any material, compound, mixture or preparation which contains any quantity of the

19 following substances having a stimulant effect on the central nervous system, including its salts,
20 isomers and salts of isomers:

21 (1) Pyrovalerone.

22 (d) Any compound, mixture or preparation containing as its single active ingredient
23 ephedrine, pseudoephedrine or phenylpropanolamine, their salts or optical isomers, or salts of
24 optical isomers except products which are for pediatric use primarily intended for administration
25 to children under the age of twelve: *Provided*, That neither the offenses set forth in section four
26 hundred one, article four of this chapter, nor the penalties therein, shall be applicable to ephedrine,
27 pseudoephedrine or phenylpropanolamine which shall be subject to the provisions of article ten
28 of this chapter.

29 (e) *Depressants*. — Unless specifically exempted or excluded or unless listed in another
30 schedule, any material, compound, mixture or preparation which contains any quantity of the
31 following substances having a depressant effect on the central nervous system, including its salts:

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

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

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

35 (4) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl] butanamide) (also referred to
36 as BRV; UCB-34714; Briviact), including its salts.

The Joint Committee on Enrolled Bills hereby certifies that the foregoing bill is correctly enrolled.

.....
Chairman, House Committee

.....
Chairman, Senate Committee

Originating in the House.

In effect ninety days from passage.

.....
Clerk of the House of Delegates

.....
Clerk of the Senate

.....
Speaker of the House of Delegates

.....
President of the Senate

The within this the.....
day of, 2017.

.....
Governor