

WEST VIRGINIA LEGISLATURE

2025 REGULAR SESSION

Introduced

House Bill 3434

By Delegate Kelly

[By Request of the Department of Homeland Security

– West Virginia State Police]

[Introduced March 17, 2025; referred to the

Committee on the Judiciary]

1 A BILL to amend and reenact §60A-2-204, §60A-2-206, §60A-2-208, §60A-2-210, and §60A-2-
2 212 of the Code of West Virginia, 1931, as amended, relating to the controlled substance
3 schedules and to clean-up errors identified in the code sections.

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 Brorphine (1-(1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2H-benzo[d]imidazol-
29 2-one);
30 Clonitazene;
31 Dextromoramide;
32 Diampromide;
33 Diethylthiambutene;
34 Difenoxin;
35 Dimenoxadol;
36 Dimepheptanol;
37 Dimethylthiambutene;
38 Dioxaphetyl butyrate;
39 Dipipanone;
40 Ethylmethylthiambutene;
41 Etonitazene;
42 Etoxidine;
43 Fentanyl analog or derivative, as that term is defined in article one of this chapter:
44 Provided, That fentanyl and carfentanil remains a Schedule II substance, as set forth in W. Va.
45 Code §60A-2-206;
46 Furethidine;

47 Hydroxypethidine;

48 Ketobemidone;

49 Levomoramide;

50 Levophenacymorphan;

51 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

52 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4-piperidiny]-phenylpropanamide);

53 Morpheridine;

54 N-Methylnorfentanyl (N-(1-Methyl-4-piperidiny)-N-phenyl-propanamide,

55 monohydrochloride);

56 ~~Norfentanyl (N-Phenyl-N-4-piperidiny-propanamide);~~

57 MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);

58 Noracymethadol;

59 Norlevorphanol;

60 Normethadone;

61 Norpipanone;

62 Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidiny] propanamide);

63 PEPAP(1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);

64 Phenadoxone;

65 Phenampromide;

66 Phenomorphan;

67 Phenoperidine;

68 Piritramide;

69 Proheptazine;

70 Properidine;

71 Propiram;

72 Racemoramide;

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

99 Thebacon.

100 (d) Hallucinogenic substances.

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

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

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

104 methoxymethamphetamine, PMMA);

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

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

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

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

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

110 acronym 25B-NBOMe;

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

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

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

114 methylphenethylamine; 2,5-DMA;

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

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

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

118 methylphenethylamine; paramethoxyamphetamine; PMA;

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

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

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

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

123 3,4-methylenedioxy amphetamine;

124 3,4-methylenedioxymethamphetamine (MDMA);

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

151 Psilocybin;

152 Psilocyn;

153 Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or
154 in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives
155 and their isomers with similar chemical structure and pharmacological activity including, but not
156 limited to the following:

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

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

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

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

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

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

164 Delta-8-tetrahydrocannabinol-O (delta-8-THC-0), Delta-9-tetrahydrocannabinol (delta-9-
165 THC-0) and Synthetic and non-naturally occurring cannabinoids.

166 The provisions of this section related to tetrahydrocannabinols are inapplicable to
167 products or substances lawfully manufactured, distributed, or possessed under the provisions of §
168 19-12E-1 et seq. and Chapter 16H of this code.

169 Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-
170 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,
171 cyclohexamine, PCE;

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

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

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

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

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

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

255 NBOMe);
256 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
257 NBOMe);
258 Synthetic Opioids (including their isomers, esters, ethers, salts and salts of isomers, esters
259 and ethers):
260 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);
261 furanyl fentanyl;
262 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
263 47700);
264 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
265 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);
266 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also known
267 as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidiny]-N-phenylpropanamide, (beta-
268 hydroxythiofentanyl);
269 N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl);
270 N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl);
271 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl);
272 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide
273 (also known as U-48800);
274 Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as
275 U-49900);
276 Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also
277 known as U-51754);
278 2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine
279 (butonitazene);
280 2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine (etodesnitazene);

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

307 JWH 018;
308 JWH 019;
309 JWH 073;
310 JWH 081;
311 JWH 122;
312 JWH 200;
313 JWH 210;
314 JWH 398;
315 AM 2201; and
316 WIN 55,212.

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

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

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

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

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

334 JWH 250;

335 JWH 203;

336 JWH 251; and

337 JWH 302.

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

341 CP 47,497 and its homologues and analogs;

342 Cannabicyclohexanol; and

343 CP 55,940.

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

348 AM 694;

349 Pravadoline WIN 48,098;

350 RCS 4; and

351 AM 679.

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

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

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

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

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

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

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

Tryptamines:

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

411 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
 412 a][1,4]diazepine);
 413 Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
 414 a][1,4]benzodiazepine);
 415 Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
 416 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
 417 a][1,4]benzodiazepine);
 418 Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-
 419 a][1,4]diazepine);
 420 Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
 421 one) ;
 422 Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); and
 423 Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
 424 a][1,4]benzodiazepine).
 425 (e) Depressants.
 426 4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-
 427 carboxamide);
 428 Alpha-Phenylacetoacetonitrile (3-Oxo-2-phenylbutanenitrile);
 429 2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-cyclohexanone,
 430 monohydrochloride);
 431 4-MEAP (2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one);
 432 Mecloqualone;
 433 Methaqualone;
 434 Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
 435 Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3
 436 a][1,4]benzodiazepine);

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

463 oxazolamine; 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine);
464 Cathinone; some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-
465 aminopropiophenone, 2-aminopropiophenone and norephedrone;
466 Ethylphenidate (ethyl 2-phenyl-2-(piperidin-2-yl)acetate);
467 Fenethylamine;
468 Mesocarb (N-phenyl-N'-(3-(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-
469 yl)carbamimidate);
470 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
471 isomers and salts of optical isomers; some other names: (2-(methylamino)-propionophenone; alpha-
472 (methylamino)propionophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-
473 methylaminopropionophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or
474 mephedrone; 3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone;
475 methylcathinone; AL-464; AL-422; AL-463 and UR1432;
476 (+-) cis-4-methylaminorex; ((+)-cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
477 N-ethylamphetamine;
478 N,N-dimethylamphetamine; also known as N,N-alpha-trimethyl-benzeneethanamine;
479 N,N-alpha-trimethylphenethylamine;
480 Alpha-pyrrolidinopropionophenone, also known as alpha-PVP, optical isomers, salts and
481 salts of isomers;
482 Substituted amphetamines:
483 2-Fluoroamphetamine;
484 3-Fluoroamphetamine;
485 4-Fluoroamphetamine;
486 2-chloroamphetamine;
487 3-chloroamphetamine;
488 4-chloroamphetamine;

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

515 phenylbutyramide);
516 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
517 yl)butyramide);
518 Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);
519 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
520 carboxamide); and
521 Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).

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

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

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

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

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

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

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

§60A-2-206.

Schedule

II.

1 (a) Schedule II consists of the drugs and other substances, by whatever official name,

common or usual name, chemical name or brand name designated, listed in this section. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including their isomers, esters, ethers, salts and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation.

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

Opium and opiate, and any salt, compound, derivative, or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrophan, nalbuphine, nalmeferene, naloxone and naltrexone, and their respective salts, but including the following:

Raw opium;

Opium extracts;

Opium fluid;

Powdered opium;

Granulated opium;

Tincture of opium;

Codeine;

Dihydroetorphine;

Ethylmorphine;

Etorphine hydrochloride;

Hydrocodone;

Hydromorphone;

Metopon;

Morphine;

28 Oripavine;

29 Oxycodone;

30 Oxymorphone; and

31 Thebaine;

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

35 Opium poppy and poppy straw;

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

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

43 (c) Opiates.

44 Alfentanil;

45 Alphaprodine;

46 Anileridine;

47 Bezitramide;

48 Bulk dextropropoxyphene (nondosage forms);

49 Carfentanil;

50 Dihydrocodeine;

51 Diphenoxylate;

52 Fentanyl;

53 Isomethadone;

54 Levo-alpha-acetylmethadol; some other names: levo-alpha-acetylmethadol, levomethadyl
55 acetate, LAAM;
56 Levomethorphan;
57 Levorphanol;
58 Metazocine;
59 Methadone;
60 Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane;
61 Moramide-Intermediate, 2-methyl-3-morpholino-1;
62 Norfentanyl;
63 Oliceridine;
64 1-diphenylpropane-carboxylic acid;
65 Pethidine; (meperidine);
66 Pethidine-Intermediate-A, 4-cyano-1-methyl-4- phenylpiperidine;
67 Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;
68 Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;
69 Phenazocine;
70 Piminodine;
71 Racemethorphan;
72 Racemorphan;
73 Remifentanil;
74 Sufentanil;
75 Tapentadol; and
76 Thiafentanil (4-(methoxycarbonyl)-4-(N-phenmethoxyacetamido)-1-2-
77 (thienyl)ethylpiperidine), including its isomers, esters, ethers, salts and salts of isomers, esters
78 and ethers.
79 (d) Stimulants.

Amphetamine, its salts, optical isomers, and salts of its optical isomers;

Methamphetamine, its salts, isomers, and salts of its isomers;

Methylphenidate;

Phenmetrazine and its salts; and

Lisdexamfetamine.

(e) Depressants.

Amobarbital;

Glutethimide;

Pentobarbital;

Phencyclidine; and

Secobarbital.

(f) Hallucinogenic substances:

Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] if in an FDA approved oral solution;

and

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

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

Immediate precursor to amphetamine and methamphetamine:

Phenylacetone;

Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;

Immediate precursors to phencyclidine (PCP):

1-phenylcyclohexylamine; and

1-piperidinocyclohexanecarbonitrile (PCC).

106 Immediate precursor to fentanyl:
107 4-anilino-N-phenethyl-4-piperidine (ANPP).

§60A-2-208.**Schedule****III.**

1 (a) Schedule III 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) Stimulants. -- Unless specifically excepted or unless listed in another schedule, any
4 material, compound, mixture or preparation which contains any quantity of the following
5 substances having a stimulant effect on the central nervous system, including its salts, isomers
6 (whether optical, position or geometric) and salts of such isomers whenever the existence of the
7 salts, isomers and salts of isomers is possible within the specific chemical designation:

8 (1) Those compounds, mixtures or preparations in dosage unit form containing any
9 stimulant substances listed in Schedule II which compounds, mixtures or preparations were listed
10 on August 25, 1971, as excepted compounds under 21 C.F.R. §1308.32, and any other drug of the
11 quantitative composition shown in that list for those drugs or which is the same except that it
12 contains a lesser quantity of controlled substances;

13 (2) Benzphetamine;

14 (3) Chlorphentermine;

15 (4) Clortermine;

16 (5) Phendimetrazine.

17 (c) Depressants. -- Unless specifically excepted or unless listed in another schedule, any
18 material, compound, mixture or preparation which contains any quantity of the following
19 substances having a depressant effect on the central nervous system:

20 (1) Any compound, mixture or preparation containing:

21 (A) Amobarbital;

22 (B) Secobarbital;

23 (C) Pentobarbital; or any salt of pentobarbital and one or more other active medicinal

- 24 ingredients which are not listed in any schedule;
- 25 (2) Any suppository dosage form containing:
- 26 (A) Amobarbital;
- 27 (B) Secobarbital;
- 28 (C) Pentobarbital; or any salt of any of these drugs and approved by the food and drug
- 29 administration for marketing only as a suppository;
- 30 (3) Any substance which contains any quantity of a derivative of barbituric acid or any salt
- 31 of barbituric acid;
- 32 (4) Aprobarbital;
- 33 (5) Butabarbital (secbutabarbital);
- 34 (6) Butalbital (including, but not limited to, Fioricet);
- 35 (7) Butobarbital (butethal);
- 36 (8) Chlorhexadol;
- 37 (9) Embutramide;
- 38 (10) Gamma Hydroxybutyric Acid preparations;
- 39 (11) Ketamine, its salts, isomers and salts of isomers [Some other names for ketamine: (+-
- 40)-2-(2-chlorophenyl)-2-(methylamino)-cyclohexanone];
- 41 (12) Lysergic acid;
- 42 (13) Lysergic acid amide;
- 43 (14) Methypylon;
- 44 (15) Perampanel, and its salts, isomers, and salts of isomers;
- 45 ~~(15)~~(16) Sulfondiethylmethane;
- 46 ~~(16)~~(17) Sulfonethylmethane;
- 47 ~~(17)~~(18) Sulfonmethane;
- 48 ~~(18)~~(19) Thiamylal;
- 49 ~~(19)~~(20) Thiopental;

~~(20)~~ (21) Tiletamine and zolazepam or any salt of tiletamine and zolazepam; some trade or other names for a tiletamine-zolazepam combination product: Telazol; some trade or other names for tiletamine: 2-(ethylamino)-2-(2-thienyl)-cyclohexanone; some trade or other names for zolazepam: 4-(2-fluorophenyl)-6, 8-dihydro-1, 3, 8-trimethylpyrazolo-[3,4-e] [1,4]-diazepin-7(1H)-one, flupyrzapon; and

~~(24)~~ (22) Vinbarbital.

(d) Nalorphine.

(e) Narcotic drugs. -- Unless specifically excepted or unless listed in another schedule:

(1) Any material, compound, mixture or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:

(A) Not more than 1.8 grams of codeine per 100 milliliters and not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium;

(B) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(C) Not more than 1.8 grams of dihydrocodeine per 100 milliliters and not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(D) Not more than 300 milligrams of ethylmorphine per 100 milliliters or not more than 15 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(E) Not more than 500 milligrams of opium per 100 milliliters or per 100 grams or not more than 25 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(F) Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams, with one

or more active, nonnarcotic ingredients in recognized therapeutic amounts.

(2) Any material, compound, mixture or preparation containing buprenorphine or its salts (including, but not limited to, Suboxone).

(f) Anabolic steroids. -- Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of anabolic steroids, including its salts, isomers and salts of isomers whenever the existence of the salts of isomers is possible within the specific chemical designation.

(g) Human growth hormones.

(h) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a United States food and drug administration approved drug product. (Some other names for dronabinol: (6aR-trans)-6a, 7, 8, 10a-tetrahydro-6, 6, 9-trimethyl-3-pentyl-6H-dibenzo [b,d] pyran-1-ol or (-)-delta-9-(trans)-tetrahydrocannabinol).

(i) Human chorionic gonadotropin, except when used for injection or implantation in cattle or any other nonhuman species and when that use is approved by the Food and Drug Administration.

§60A-2-210.

Schedule

IV.

(a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including their isomers, esters, ethers, salts and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation.

(b) Narcotic drugs. — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:

11 Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate
12 per dosage unit; and

13 Dextropropoxyphene (alpha-()-4-dimethylamino-1,2-diphenyl-3-methyl-2-
14 propionoxybutane).

15 (c) Depressants.

16 Alfaxalone;

17 Alprazolam;

18 Barbitol;

19 Bromazepam;

20 Camazepam;

21 Carisoprodol;

22 Chloral betaine;

23 Chloral hydrate;

24 Chlordiazepoxide;

25 Clobazam;

26 Clonazepam;

27 Clorazepate;

28 Clotiazepam;

29 Cloxazolam;

30 Daridorexant;

31 Delorazepam;

32 Diazepam;

33 Dichloralphenazone;

34 Estazolam;

35 Ethchlorvynol;

36 Ethinamate;

- 37 Ethyl loflazepate;
- 38 Fludiazepam;
- 39 Flunitrazepam;
- 40 Flurazepam;
- 41 Fospropofol;
- 42 Halazepam;
- 43 Haloxazolam;
- 44 Ketazolam;
- 45 Lemborexant.
- 46 Loprazolam;
- 47 Lorazepam;
- 48 Lormetazepam;
- 49 Mebutamate;
- 50 Medazepam;
- 51 Meprobamate;
- 52 Methohexital;
- 53 Methylphenobarbital (mephobarbital);
- 54 Midazolam;
- 55 Nimetazepam;
- 56 Nitrazepam;
- 57 Nordiazepam;
- 58 Oxazepam;
- 59 Oxazolam;
- 60 Paraldehyde;
- 61 Petrichloral;
- 62 Phenobarbital;

- 63 Pinazepam;
- 64 Prazepam;
- 65 Quazepam;
- 66 Remimazolam.
- 67 Temazepam;
- 68 Tetrazepam;
- 69 Triazolam;
- 70 Xylazine;
- 71 Zaleplon;
- 72 Zolpidem;
- 73 Zopiclone; and
- 74 Suvorexant ([[(7R)-4-(5-chloro-1,3-benzoxazol-2-yl)-7-methyl-1,4-diazepan-1-yl] [5-
- 75 methyl-2-(2H-1,2,3-triazol-2-yl)phenyl]methanone).
- 76 Zuranolone;
- 77 (d) Any material, compound, mixture, or preparation which contains any quantity of
- 78 Fenfluramine and Dexfenfluramine.
- 79 (e) Stimulants.
- 80 Cathine (()-norpseudoephedrine);
- 81 Diethylpropion;
- 82 Fencamfamin;
- 83 Fenproporex;
- 84 Mazindol;
- 85 Mefenorex;
- 86 Modafinil;
- 87 Pemoline (including organometallic complexes and chelates thereof);
- 88 Phentermine;

Pipradrol;

Serdexmethylphenidate;

Sibutramine;

SPA ((-)-1-dimethylamino-1,2-diphenylethane); and

Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-oxopropyl

[(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid);

(f) Other substances.

Lorcaserin;

Pentazocine;

Butorphanol;

Tramadol (2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol); and

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

§60A-2-212.

Schedule

V.

(a) Schedule V shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation which contains any quantity of the following substances, including their isomers, esters, ethers, salts and salts of isomers, esters and ethers, whenever the existence of such isomers, esters, ethers and salts is possible within the specific chemical designation.

(b) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture or preparation containing any of the following narcotic drugs or their salts calculated as the free anhydrous base or alkaloid in limited quantities as set forth below, which shall include one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the

compound, mixture or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone.

Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;

Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams;

Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams;

Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit;

Not more than 100 milligrams of opium per 100 milliliters or per 100 grams; and

Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.

(c) Stimulants:

Pyrovalerone.

(d) Any compound, mixture, or preparation containing as its single active ingredient ephedrine, pseudoephedrine, or phenylpropanolamine, their salts or optical isomers, or salts of optical isomers except products which are for pediatric use primarily intended for administration to children under the age of 12: *Provided*, That neither the offenses set forth in section four hundred one, article four of this chapter, nor the penalties therein, shall be applicable to ephedrine, pseudoephedrine or phenylpropanolamine which shall be subject to the provisions of article ten of this chapter.

(e) Depressants:

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

Ganaxolone (3 α -hydroxy-3 β -methyl-5 α -pregnan-20-one);

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

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

(f) Other substances:

- 37 Gabapentin;
- 38 Pregabalin;
- 39 Cenobamate; and
- 40 Lasmiditan.

NOTE: The purpose of this bill is to clean-up errors identified in the code sections.

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