WEST VIRGINIA LEGISLATURE

2020 REGULAR SESSION

Introduced

House Bill 4858

By Delegates Steele, Higginbotham, Dean, Hill,
Pack, Toney, McGeehan, Storch, Linville, Hamrick
AND Maynard

[Introduced February 11, 2020; Referred to the Committee on Health and Human Resources then the Judiciary]

A BILL to amend and reenact §60A-2-204 and §60A-2-210 of the Code of West Virginia, 1931, as amended, all relating to classifying "marihuana" and tetrahydrocannabinols as a Schedule IV controlled substance; deleting marihuana and tetrahydrocannabinols from Schedule I listing; and adding these substances to the Schedule IV list under other substances.

Be it enacted by the Legislature of West Virginia:

ARTICLE 2. STANDARDS AND SCHEDULES.

§60A-2-204. Schedule I.

18

Benzethidine;

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 existence 4 of such isomers, esters, ethers and salts is possible within the specific chemical designation. 5 (b) Opiates. 6 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl) -4-piperidinyl]— 7 phenylacetamide); 8 Acetylmethadol; 9 Allylprodine: 10 Alphacetylmethadol (except levoalphacetylmethadol also known as levo-alpha-11 acetylmethadol, levomethadyl acetate, or LAAM); 12 Alphameprodine; 13 Alphamethadol; 14 Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propionanilide: 15 1-(1-methyl-2-phenylethyl)-4-((propanilido) piperidine); 16 (N-[1-methyl-2-(2-thienyl) Alpha-methylthiofentanyl ethyl-4-piperidinyl]— 17 phenylpropanamide);

19	Betacetylmethadol;
20	Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl) -4- piperidinyl]-N-
21	phenylpropanamide);
22	Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2- hydroxy-2-phenethyl)-3-methyl-4-
23	piperidinyl]-N-phenylpropanamide);
24	Betameprodine;
25	Betamethadol;
26	Betaprodine;
27	Clonitazene;
28	Dextromoramide;
29	Diampromide;
30	Diethylthiambutene;
31	Difenoxin;
32	Dimenoxadol;
33	Dimepheptanol;
34	Dimethylthiambutene;
35	Dioxaphetyl butyrate;
36	Dipipanone;
37	Ethylmethylthiambutene;
38	Etonitazene;
39	Etoxeridine;
40	Furethidine;
41	Hydroxypethidine;
42	Ketobemidone;
43	Levomoramide;
44	Levophenacylmorphan;

45	3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4- piperidyl]-N-phenylpropanamide);
46	3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4- piperidinyl]—phenylpropanamide);
47	Morpheridine;
48	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
49	Noracymethadol;
50	Norlevorphanol;
51	Normethadone;
52	Norpipanone;
53	Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2- phenethyl)-4-piperidinyl] propanamide);
54	PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
55	Phenadoxone;
56	Phenampromide;
57	Phenomorphan;
58	Phenoperidine;
59	Piritramide;
60	Proheptazine;
61	Properidine;
62	Propiram;
63	Racemoramide;
64	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4- piperidinyl]-propanamide);
65	Tilidine;
66	Trimeperidine.
67	(c) Opium derivatives:
68	Acetorphine;
69	Acetyldihydrocodeine;
70	Benzylmorphine;

71	Codeine methylbromide;
72	Codeine-N-Oxide;
73	Cyprenorphine;
74	Desomorphine;
75	Dihydromorphine;
76	Drotebanol;
77	Etorphine (except HCl Salt);
78	Heroin;
79	Hydromorphinol;
80	Methyldesorphine;
81	Methyldihydromorphine;
82	Morphine methylbromide;
83	Morphine methylsulfonate;
84	Morphine-N-Oxide;
85	Myrophine;
86	Nicocodeine;
87	Nicomorphine;
88	Normorphine;
89	Pholcodine;
90	Thebacon.
91	(d) Hallucinogenic substances.
92	Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-
93	indole-3-ethanamine; 3-(2- aminobutyl) indole; alpha-ET; and AET;
94	4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
95	dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;
96	4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-

97 dimethoxyphenyl)-1-aminoethane: alpha- desmethyl DOB: 2C-B. Nexus: 98 N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the 99 acronym 25B-NBOMe. 100 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe) 101 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe) 102 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-103 methylphenethylamine; 2,5-DMA; 104 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET; 105 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7); 106 4-methoxyamphetamine: some trade or other names: 4-methoxy-alpha-107 methylphenethylamine; paramethoxyamphetamine; PMA; 108 5-methoxy-3, 4-methylenedioxy-amphetamine; 109 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-110 dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP"; 111 3,4-methylenedioxy amphetamine; 112 3,4-methylenedioxymethamphetamine (MDMA); 113 3,4-methylenedioxy-N-ethylamphetamine (also known as (ethyl-alpha-methyl-3,4 114 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA); 115 N-hydroxy-3,4-methylenedioxyamphetamine (also known as (hydroxy-alpha-methyl-3,4 116 (methylenedioxy) phenethylamine, and (hydroxy MDA); 117 3.4.5-trimethoxy amphetamine: 118 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT): 119 Alpha-methyltryptamine (other name: AMT); 120 Bufotenine; some trade other 3-(beta-Dimethylaminoethyl)-5and names: 121 hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-122 dimethyltryptamine; mappine;

123	Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;
124	Dimethyltryptamine; some trade or other names: DMT;
125	5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);
126	Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-
127	methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;
128	Lysergic acid diethylamide;
129	Marihuana;
130	Mescaline;
131	Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,
132	6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;
133	Peyote; meaning all parts of the plant presently classified botanically as Lophophora
134	williamsii Lemaire, whether growing or not, the seeds thereof, any extract from any part of such
135	plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of
136	such plant, its seeds or extracts;
137	N-ethyl-3-piperidyl benzilate;
138	N-methyl-3-piperidyl benzilate;
139	Psilocybin;
140	Psilocyn;
141	Tetrahydrocannabinols; synthetic Synthetic equivalents of the substances contained in the
142	plant, or in the resinous extractives of Cannabis sp. and/or synthetic substances, immediate
143	derivatives and their isomers with similar chemical structure and pharmacological activity such as
144	the following:
145	delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;
146	delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;
147	delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;
148	(Since nomenclature of these substances is not internationally standardized, compounds

149 of these structures, regardless of numerical designation of atomic positions covered.) Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-150 151 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, 152 cyclohexamine, PCE; 153 Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-154 pyrrolidine, PCPv, PHP; 155 Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-156 cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine; TPCP, TCP; 157 1[1-(2-thienyl)cyclohexyl]pyrroldine; some other names: TCPy. 158 4-methylmethcathinone (Mephedrone): 159 3.4-methylenedioxypyrovalerone (MDPV); 160 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E); 161 2-(2.5-Dimethoxy-4-methylphenyl)ethanamine (2C-D) 162 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C) 163 2-(4-lodo-2,5-dimethoxyphenyl)ethanamine (2C-I) 164 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2) 165 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4) 166 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H) 167 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N) 168 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P) 169 3,4-Methylenedioxy-N-methylcathinone (Methylone) 170 2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, itsoptical isomers, salts and 171 salts of isomers 172 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-173 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT) 174 Alpha-methyltryptamine (other name: AMT)

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175
              5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT)
              Synthetic Cannabinoids as follows:
176
177
              2-[(1R,3S)-3-hydroxycyclohexyl]-5- (2-methyloctan-2-yl)phenol) {also known as CP
178
       47,497 and homologues);
179
              rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl)phenol {also known as CP
180
       47,497-C8 homolog);
181
              [(6aR)-9-(hydroxymethyl)-6,
                                               6-dimethyl-3-(2-methyloctan-2-yl)-6a,
                                                                                         7,10,10a-
182
       tetrahydrobenzo[c]chromen-1-ol)] {also known as HU-210};
183
              (dexanabinol);
184
              (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
185
       tetrahydrobenzol[c]chromen-1-ol) {also known as HU-211};
186
              1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};
187
              1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};
188
              (2-methyl-1-propyl-1H-indol-3-yl)-1-napthalenyl-methanone {also known as JWH-015};
189
              (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};
190
              [1-[2-(4-morpholinyl) ethyl] -1H-indol-3-yl]-1-naphthalenyl-methanone {also known as
191
       JWH-200};
192
              1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};
193
              2-((1S,2S,5S)-5-hydroxy-2- (3-hydroxtpropyl)cyclohexyl) -5-(2-methyloctan-2-yl)phenol
194
       {also known as CP 55,940};
195
              (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-122};
196
              (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-398;
197
              (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};
198
              1-(1-(2-cyclohexylethyl) -1H-indol-3-yl) -2-(2-methoxyphenyl) ethanone {also known as
199
       RCS-8};
200
              1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);
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201	1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and
202	1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).
203	Synthetic cannabinoids:
204	CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
205	YL)phenol);
206	HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,
207	10A-tetrahydrobenzo[C] chromen-1-OL)];
208	HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-
209	YL)-6A,7,10,10atetrahydrobenzo[C]chromen-1-OL);
210	JWH-018, 1-pentyl-3-(1-naphthoyl)indole;
211	JWH-019, 1-hexyl-3-(1-naphthoyl)indole;
212	JWH-073, 1-butyl-3-(1-naphthoyl)indole;
213	JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)- Naphthalen-1-ylmethanone;
214	JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.]
215	Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-
216	ADB);
217	Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB);
218	Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-
219	AMB);
220	N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA);
221	N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
222	(ADB-FUBINACA);
223	Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
224	(MDMB-CHMICA);
225	Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate
226	(MDMB-FUBINACA);

227	Tetrahydrocannabinols:
228	DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.
229	DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.
230	DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.
231	Synthetic Phenethylamines
232	2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
233	NBOMe);
234	2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-
235	NBOMe);
236	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
237	NBOMe);
238	Synthetic Opioids (icluding their isomers, esters, ethers, salts and salts of isomers, esters
239	and ethers):
240	N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);
241	furanyl fentanyl;
242	3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
243	47700);
244	N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
245	phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);
246	N-[1-[2-hydroxy-2-(thiophen-2-yl)ethylpiperidin-4-yl]-N-phenylpropionamide, also known
247	as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-
248	hydroxythiofentanyl).
249	N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl)
250	N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl)
251	N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl)
252	2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide

253 (also known as U-48800) 254 Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as 255 U-49900) 256 Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also 257 known as U-51754) 258 Opioid Receptor Agonist 259 AH-7921 (3,4-dichloro-N- (1dimethylamino)cyclohexylmethyl]benzamide). 260 Naphthoylindoles or any compound containing a 3-(-1- Napthoyl) indole structure with 261 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole 262 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall 263 include the following: 264 JWH 015; 265 JWH 018; 266 JWH 019; 267 JWH 073; 268 JWH 081; 269 JWH 122; 270 JWH 200; 271 JWH 210; 272 JWH 398; 273 AM 2201; 274 WIN 55,212. 275 Naphylmethylindoles or any compound containing a 1hindol-3-yl-(1-naphthyl) methane 276 structure with a substitution at the nitrogen atom of the indole ring whether or not further 277 substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to 278 any extent. This shall include, but not be limited to, JWH 175 and JWH 184.

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

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

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

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

292 JWH 250;

293 JWH 203;

294 JWH 251;

295 JWH 302.

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

CP 47,497 and its homologues and analogs;

Cannabicyclohexanol;

301 CP 55,940.

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

305	the following:
306	AM 694;
307	Pravadoline WIN 48,098;
308	RCS 4;
309	AM 679.
310	[2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-
311	napthalenymethanone. This shall include WIN 55,212-2.
312	Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol
313	structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051
314	and JWH 133.
315	Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with
316	substitution at the nitrogen atom of the indole ring whether or not further substituted in the
317	adamantoyl ring system to any extent. This shall include AM1248.
318	Tetramethylcyclopropylindoles or any compound containing A 3-
319	tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring
320	whether or not further substituted in the indole ring to any extent and whether or not substituted
321	in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.
322	N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.
323	Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as
324	demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and
325	V, not federal Food and Drug Administration approved drug or used within legitimate, approved
326	medical research. Since nomenclature of these substances is not internationally standardized,
327	any immediate precursor or immediate derivative of these substances shall be covered.
328	Tryptamines:
329	5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)
330	4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT)

331	4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)
332	4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)
333	4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT)
334	5-methoxy-α-methyltryptamine (5-MeO-AMT)
335	4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT)
336	4-hydroxy Diethyltryptamine (4-HO-DET)
337	5- methoxy- N,N- diallyltryptamine (5-MeO-DALT)
338	4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT)
339	4-hydroxy Diethyltryptamine (4-HO-DET)
340	(e) Depressants.
341	Mecloqualone;
342	Methaqualone.
343	(f) Stimulants.
344	Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
345	dihydro-5-phenyl-2-oxazolamine;
346	Cathinone; some trade or other names: 2-amino-1-phenyl-1- propanone, alpha-
347	aminopropiophenone, 2-aminopropiophenone and norephedrone;
348	Fenethylline;
349	Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
350	isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-
351	(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—-
352	methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or
353	mephedrone;3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone;
354	methylcathinone; AL-464; AL-422; AL- 463 and UR1432;
355	(+-) cis-4-methylaminorex; ((+-)cis-4,5-dihydro-4-methyl- 5-phenyl-2-oxazolamine);
356	N-ethylamphetamine;

357 N,N-dimethylamphetemine; also known as N,N-alpha- trimethyl-benzeneethanamine; 358 N,N-alpha-trimethylphenethylamine. 359 Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and 360 salts of isomers. 361 Substituted amphetamines: 362 2-Fluoroamphetamine 363 3-Fluoroamphetamine 364 4-Fluoroamphetamine 365 2-chloroamphetamine 366 3-chloroamphetamine 367 4-chloroamphetamine 368 2-Fluoromethamphetamine 369 3-Fluoromethamphetamine 370 4-Fluoromethamphetamine 371 4-chloromethamphetamine 372 (g) Temporary listing of substances subject to emergency scheduling. Any material, 373 compound, mixture or preparation which contains any quantity of the following substances: N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts, 374 375 and salts of isomers. 376 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical 377 isomers, salts and salts of isomers. 378 N-benzylpiperazine, also known as BZP. 379 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide); 380 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-381 butyramide); 382 Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);

383	Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-
384	acetamide);
385	3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-
386	phenylbutyramide);
387	4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
388	yl)butyramide);
389	Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);
390	Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
391	carboxamide);
392	Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).
393	(h) The following controlled substances are included in Schedule I:
394	Synthetic Cathinones or any compound, except bupropion or compounds listed under a
395	different schedule, or compounds used within legitimate and approved medical research,
396	structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic
397	or fused polycyclic ring systems, whether or not the compound is further modified in any of the
398	following ways:
399	By substitution in the ring system to any extent with Alkyl, alkylenedioxy, alkoxy, haloalkyl,
400	hydroxyl or halide Substituents whether or not further substituted in the ring system by one or
401	more other univalent substituents.
102	By substitution at the 3-position with an acyclic alkyl substituent.
403	By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl or methoxybenzyl
104	groups.
405	By inclusion of the 2-amino nitrogen atom in a cyclic structure.
406	Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as
107	demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and
108	V, not federal Food and Drug Administration approved drug or used within legitimate, approved

409 medical research.

24

Clobazam;

§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. 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. — Unless specifically excepted or unless listed in another schedule, 8 any material, compound, mixture or preparation containing any of the following narcotic drugs, or 9 their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth 10 below: 11 Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate 12 per dosage unit; 13 (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-Dextropropoxyphene 14 propionoxybutane). 15 (c) Depressants. 16 Alprazolam: 17 Barbital; 18 Bromazepam; 19 Camazepam; 20 Carisoprodol: 21 Chloral betaine; 22 Chloral hydrate; 23 Chlordiazepoxide;

25	Clonazepam;
26	Clorazepate;
27	Clotiazepam;
28	Cloxazolam;
29	Delorazepam;
30	Diazepam;
31	Dichloralphenazone;
32	Estazolam;
33	Ethchlorvynol;
34	Ethinamate;
35	Ethyl loflazepate;
36	Fludiazepam;
37	Flunitrazepam;
38	Flurazepam;
39	Fospropofol;
40	Halazepam;
41	Haloxazolam;
42	Ketazolam;
43	Loprazolam;
44	Lorazepam;
45	Lormetazepam;
46	Mebutamate;
47	Medazepam;
48	Meprobamate;
49	Methohexital;
50	Methylphenobarbital (mephobarbital);

51	Midazolam;
52	Nimetazepam;
53	Nitrazepam;
54	Nordiazepam;
55	Oxazepam;
56	Oxazolam;
57	Paraldehyde;
58	Petrichloral;
59	Phenobarbital;
60	Pinazepam;
61	Prazepam;
62	Quazepam;
63	Temazepam;
64	Tetrazepam;
65	Triazolam;
66	Zaleplon;
67	Zolpidem;
68	Zopiclone'
69	Suvorexant ([(7R)-4-(5-chloro-1,3-benzoxazol-2-yl)-7-methyl-1,4-diazepan-1-yl] [5-
70	methyl-2-(2H-1,2,3-triazol-2-yl)phenyl]methanone).
71	(d) Any material, compound, mixture or preparation which contains any quantity of
72	Fenfluramine and Dexfenfluramine.
73	(e) Stimulants.
74	Cathine ((+)-norpseudoephedrine);
75	Diethylpropion;
76	Fencamfamin;

77	Fenproporex;
78	Mazindol;
79	Mefenorex;
80	Modafinil;
81	Pemoline (including organometallic complexes and chelates thereof);
82	Phentermine;
83	Pipradrol;
84	Sibutramine;
85	SPA ((-)-1-dimethylamino-1,2-diphenylethane);
86	Eluxadoline (5-[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropy
87	[(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid);
88	(f) Other substances. —
89	Pentazocine;
90	Butorphanol;
91	Tramadol (2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol);
92	Amyl nitrite, butyl nitrite, isobutyl nitrite and the other organic nitrites are controlled
93	substances and no product containing these compounds as a significant component shall be
94	possessed, bought or sold other than pursuant to a bona fide prescription or for industrial or
95	manufacturing purposes;
96	Marijuana and plants, or parts of plants, of the genus Cannabis; and
97	Tetrahydrocannabinols.

NOTE: The purpose of this bill is to reschedule marijuana (marihuana) as a Schedule IV rather than a Schedule I controlled substance and specify it as a substance that is not a narcotic.

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