

# **WEST VIRGINIA LEGISLATURE**

## **2026 REGULAR SESSION**

**Introduced**

### **House Bill 5146**

By Delegate Young

[Introduced February 03, 2026; referred to the  
Committee on Health and Human Resources then the  
Judiciary]

A BILL to amend and reenact §60A-2-204 of the Code of West Virginia, 1931, as amended, relating to removing certain substances from schedule I of the Uniform Controlled Substances Act, including marihuana, Psilocybin, and Tetrahydrocannabinols.

*Be it enacted by the Legislature of West Virginia:*

**ARTICLE 2. STANDARDS AND SCHEDULES.**

<b>§60A-2-204.</b>	<b>Schedule</b>	<b>I.</b>
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(a) Schedule I 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 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) Opiates.

Acetyl-alpha-methylfentanyl(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-phenylacetamide);

Acetylmethadol;

Allylprodine;

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

Alphameprodine;

Alphamethadol;

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

Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-phenylpropanamide);

Benzethidine;

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           Borophine (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  
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-methoxymethamphetamine,

104 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; some trade 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 Paraheptyl-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;

Psilocybin;

Psilocyn;

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

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

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

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

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

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

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

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

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

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

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

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

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



177 4-methylmethcathinone (Mephedrone);  
178 3,4-methylenedioxypyrovalerone (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-Napthoyl) 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 Naphylmethyloindoles 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 Naphthylmethyloindenes 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- $\alpha$ -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 Diclazepam (7-Chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-  
456 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-oxazamine; 4-

463 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-yl)carbamimidate);  
469 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical  
470 isomers and salts of optical isomers; some other names: (2-(methylamino)-propionophenone; alpha-  
471 (methylamino)propionophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-  
472 methylaminopropionophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or  
473 mephedrone; 3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone;  
474 methylcathinone; AL-464; AL-422; AL-463 and UR1432;  
475 (+-) cis-4-methylaminorex; ((+)-)cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);  
476 N-ethylamphetamine;  
477 N,N-dimethylamphetamine; also known as N,N-alpha-trimethyl-benzeneethanamine;  
478 N,N-alpha-trimethylphenethylamine;  
479 Alpha-pyrrolidinopropionophenone, also known as alpha-PVP, optical isomers, salts and  
480 salts of isomers;  
481 Substituted amphetamines:  
482 2-Fluoroamphetamine;  
483 3-Fluoroamphetamine;  
484 4-Fluoroamphetamine;  
485 2-chloroamphetamine;  
486 3-chloroamphetamine;  
487 4-chloroamphetamine;  
488 2-Fluoromethamphetamine;

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

515 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-  
516 yl)butyramide);

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

518 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-  
519 carboxamide); and

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

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

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

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

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

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

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

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

NOTE: The purpose of this bill is to remove certain substances from schedule I of the Uniform Controlled Substances Act, including marihuana, Psilocybin, and Tetrahydrocannabinols.

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