# **WEST VIRGINIA LEGISLATURE**

# **2022 REGULAR SESSION**

## Introduced

# **Senate Bill 703**

By Senators Trump, Weld, and Phillips

[Introduced February 21, 2022; referred

to the Committee on Health and Human Resources]

1 A BILL to amend and reenact \$60A-2-204: 60A-2-206: 60A-2-210. and 60A-2-212 of the Code 2 of West Virginia, 1931, as amended, all relating to controlled substances, adding the 3 following to the schedule I substance list: N-Methylnorfentanyl(N-(1-Methyl-4-piperidinyl)-4 N-phenyl-propanamide, monohydrochloride); Norfentanyl (N-Phenyl-N-4-piperidinyl-5 propanamide); 3-Hydoxy-phencyclidine (other name hydroxy PCP); Marijuana (Cannabis, 6 sp.); FDU-PB-22 (1-Naphthyl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate); FUB-PB-22 7 1-(4-fluorobenzyl)-1H-indole-3-carboxylate); 5-Fluoro-MN-24 (Quinolin-8-yl (1-(5-8 Fluoropentyl)-N-(naphthalen-1-yl)-1H-indole-3-carboxamide); MN-24 (N-(naphthalen-1-9 yl)-1-pentyl-1H-indole-3-carboxamide); SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-10 3-carboxylate); SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide); Methyl-11 Ethylaminopentiophenone; **FUB-AMB** (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-12 carbonyl)-L-valinate); 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-5F-AB-PINACA 13 indole-3-carboxylate); (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-14 fluoropentyl)-1H-indazole-3carboxamide); MMB-CHMICA (Methyl 2-(1-15 (cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoat); MN-24 (N-(naphthalen-16 1-vI)-1-pentyl-1H-indole-3-carboxamide): **SDB-005** (Naphthalen-1-vl 1-pentyl-1H-17 indazole-3-carboxylate); SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide); (2-(ethylamino)-1-phenyl-1-propanone, 18 Ethcathinone monohydrochloride); Methyl-19 Ethylaminopentiophenone; **FUB-AMB** (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-20 carbonyl)-L-valinate); 5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-21 indole-3-carboxylate);5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-22 fluoropentyl)-1H-indazole-3carboxamide); MMB-CHMICA (Methyl 2-(1-23 (cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoat); Bromazolam (8-bromo-24 1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Clonazolam (6-(2-25 chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); 26 Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-

27	benzodiazepin-2-one); Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f]
28	[1,2,4]triazolo[4,3-a][1,4]diazepine); Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-
29	4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Flubromazepam (7-bromo-5-(2-
30	fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one); Flubromazolam (8-bromo-6-(2-
31	fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Flunitrazolam (6-(2-
32	fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-a][1,4]diazepine); Nifoxipam
33	(5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-one);
34	Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
35	Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
36	a][1,4]benzodiazepine); 4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-
37	2- yl)-1H-indazole-3-carboxamide); Alpha-Phenylacetoacetonitrile (3-Oxo-2-
38	phenylbutanenitrile); 2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-
39	cyclohexanone, monohydrochloride); 4-MEAP (2-(Ethylamino)-1-(4-
40	methylphenyl)pentan-1-one); Bromazolam (8-bromo-1-methyl-6-phenyl-4H-
41	[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Clonazolam (6-(2-chlorophenyl)-1-methyl-8-
42	nitro-4 H-[1,2,4]triazolo[4,3 a][1,4]benzodiazepine); Cloniprazepam (5-(2-chlorophenyl)-1-
43	(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-benzodiazepin-2-one); Etizolam (4-(2-
44	chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-a][1,4]diazepine);
45	Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
46	a][1,4]benzodiazepine); Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-
47	1,4-benzodiazepin-2-one); Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-
48	[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-
49	nitro-4H-benzo[f][1,2,4]triazolo[4,3-a][1,4]diazepine); Nifoxipam (5-(2-fluorophenyl)-1,3-
50	dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-one); Nitrazolam (1-methyl-8-nitro-6-
51	phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Pyrazolam (8-bromo-1-methyl-6-(2-
52	pyridinyl)-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine); Declazepam (7-Chloro-5-(2-

53 chlorophenyl)-1-methyl-1.3-dihydro-2H-1.4-benzodiazepin-2-one); Deschloroetizolam (2-54 Ethyl-9-methyl-4-phenyl-6H-thieno[3,2f][1,2,4]triazolo[4,3-a][1,4]diazepine); 55 Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride); Alpha-PHP (1-MPHP 56 Phenyl-2-(pyrrolidin-1-yl)hexan-1-one); (1-(4-Methylphenyl)-2-(pyrrolidin-1-57 yl)hexan-1-one); PV8 (1-Phenyl-2-(pyrrolidin-1-yl)heptan-1-one); 4-Chloro-Alpha-PVP (1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one); N-Ethylhexedrone (2-(Ethylamino)-1-58 59 phenylhexan-1-one); Methoxetamine (2-(Ethylamino)-2-(3-methoxyphenyl)-60 cyclohexanone); 3-Fluorophenmetrazine (2-(3-Fluorophenyl)-3-methylmorpholine); 61 adding the following to the schedule II list: Norfentanyl; Oliceridine; adding the following 62 to the schedule IV list: Lemborexant; Remimazolam; Serdexmethylphenidate; and 63 removing Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid] from the schedule V 64 list.

Be it enacted by the Legislature of West Virginia:

#### ARTICLE 2. STANDARDS AND SCHEDULES.

#### §60A-2-204. Schedule I.

- (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.
- 5 (b) Opiates.

1

2

3

4

- 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	Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl) ethyl-4-piperidinyl]—
17	phenylpropanamide);
18	Benzethidine;
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	N-Methylnorfentanyl (N-(1-Methyl-4-piperidinyl)-N-phenyl-propanamide,
49	monohydrochloride);
50	Norfentanyl (N-Phenyl-N-4-piperidinyl-propanamide);
51	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
52	Noracymethadol;
53	Norlevorphanol;
54	Normethadone;
55	Norpipanone;
56	Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2- phenethyl)-4-piperidinyl] propanamide);
57	PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
58	Phenadoxone;
59	Phenampromide;
60	Phenomorphan;
61	Phenoperidine;
62	Piritramide;
63	Proheptazine;

64	Properidine;
65	Propiram;
66	Racemoramide;
67	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4- piperidinyl]-propanamide);
68	Tilidine;
69	Trimeperidine.
70	(c) Opium derivatives:
71	Acetorphine;
72	Acetyldihydrocodeine;
73	Benzylmorphine;
74	Codeine methylbromide;
75	Codeine-N-Oxide;
76	Cyprenorphine;
77	Desomorphine;
78	Dihydromorphine;
79	Drotebanol;
80	Etorphine (except HCl Salt);
81	Heroin;
82	Hydromorphinol;
83	Methyldesorphine;
84	Methyldihydromorphine;
85	Morphine methylbromide;
86	Morphine methylsulfonate;
87	Morphine-N-Oxide;
88	Myrophine;
89	Nicocodeine;

90	Nicomorphine;
91	Normorphine;
92	Pholcodine;
93	Thebacon.
94	(d) Hallucinogenic substances.
95	Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-
96	indole-3-ethanamine; 3-(2- aminobutyl) indole; alpha-ET; and AET;
97	4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
98	dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;
99	4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-
100	dimethoxyphenyl)-1-aminoethane; alpha- desmethyl DOB; 2C-B, Nexus;
101	N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the
102	acronym 25B-NBOMe.
103	2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe)
104	2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)
105	2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-
106	methylphenethylamine; 2,5-DMA;
107	2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;
108	2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
109	4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-
110	methylphenethylamine; paramethoxyamphetamine; PMA;
111	3-Hydoxy-phencyclidine (other name hydroxy PCP)
112	5-methoxy-3, 4-methylenedioxy-amphetamine;
113	4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-
114	dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";
115	3,4-methylenedioxy amphetamine;

116	3,4-methylenedioxymethamphetamine (MDMA);
117	3,4-methylenedioxy-N-ethylamphetamine (also known as ( ethyl-alpha-methyl-3,4
118	(methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);
119	N-hydroxy-3,4-methylenedioxyamphetamine (also known as ( hydroxy-alpha-methyl-3,4
120	(methylenedioxy) phenethylamine, and ( hydroxy MDA);
121	3,4,5-trimethoxy amphetamine;
122	5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
123	Alpha-methyltryptamine (other name: AMT);
124	Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-
125	hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-
126	dimethyltryptamine; mappine;
127	Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;
128	Dimethyltryptamine; some trade or other names: DMT;
129	5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);
130	Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-
131	methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;
132	Lysergic acid diethylamide;
133	Marihuana; Marijuana (Cannabis, sp.);
134	Mescaline;
135	Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,
136	6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;
137	Peyote; meaning all parts of the plant presently classified botanically as Lophophora
138	williamsii Lemaire, whether growing or not, the seeds thereof, any extract from any part of such
139	plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of
140	such plant, its seeds or extracts;
141	N-ethyl-3-piperidyl benzilate;

142	N-methyl-3-piperidyl benzilate;
143	Psilocybin;
144	Psilocyn;
145	Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or
146	in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives
147	and their isomers with similar chemical structure and pharmacological activity such as the
148	following:
149	delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;
150	delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;
151	delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;
152	(Since nomenclature of these substances is not internationally standardized, compounds
153	of these structures, regardless of numerical designation of atomic positions covered.)
154	Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-
155	phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,
156	cyclohexamine, PCE;
157	Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-
158	pyrrolidine, PCPy, PHP;
159	Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-
160	cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine; TPCP, TCP;
161	1[1-(2-thienyl)cyclohexyl]pyrroldine; some other names: TCPy.
162	4-methylmethcathinone (Mephedrone);
163	3,4-methylenedioxypyrovalerone (MDPV);
164	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);
165	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D)
166	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C)
167	2-(4-lodo-2,5-dimethoxyphenyl)ethanamine (2C-l)

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168
              2-[4-(Ethylthio)-2.5-dimethoxyphenyllethanamine (2C-T-2)
169
              2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4)
170
              2-(2,5-Dimethoxyphenyl)ethanamine (2C-H)
171
              2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N)
172
              2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P)
173
              3,4-Methylenedioxy-N-methylcathinone (Methylone)
174
              2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, itsoptical isomers, salts and
175
       salts of isomers
176
              5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-
177
       (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT)
178
              Alpha-methyltryptamine (other name: AMT)
179
              5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT)
180
              Synthetic Cannabinoids as follows:
181
              2-[(1R,3S)-3-hydroxycyclohexyl]-5- (2-methyloctan-2-yl)phenol) {also known as CP
182
       47,497 and homologues);
183
              rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl)phenol {also known as CP
184
       47,497-C8 homolog);
185
                                                                                        7,10,10a-
             [(6aR)-9-(hydroxymethyl)-6,
                                              6-dimethyl-3-(2-methyloctan-2-yl)-6a,
186
       tetrahydrobenzo[c]chromen-1-ol)] {also known as HU-210};
187
              (dexanabinol);
188
              (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
189
       tetrahydrobenzol[c]chromen-1-ol) {also known as HU-211};
190
              1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};
191
              1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};
192
             (2-methyl-1-propyl-1H-indol-3-yl)-1-napthalenyl-methanone {also known as JWH-015};
193
             (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};
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194
              [1-[2-(4-morpholinyl) ethyl] -1H-indol-3-yl]-1-naphthalenyl-methanone {also known as
195
       JWH-200};
196
              1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};
197
              2-((1S,2S,5S)-5-hydroxy-2- (3-hydroxtpropyl)cyclohexyl) -5-(2-methyloctan-2-yl)phenol
198
       {also known as CP 55.940}:
199
              (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-122};
200
              (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-398;
201
              (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};
202
              1-(1-(2-cyclohexylethyl) -1H-indol-3-yl) -2-(2-methoxyphenyl) ethanone {also known as
203
       RCS-8};
204
              1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);
205
              1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and
206
              1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).
207
              Synthetic cannabinoids:
208
              CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-
209
              YL)phenol);
210
              HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,
211
       10A-tetrahydrobenzo[C] chromen-1-OL)]:
212
              HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-
213
       YL)-6A,7,10,10atetrahydrobenzo[C]chromen-1-OL);
214
              JWH-018, 1-pentyl-3-(1-naphthoyl)indole;
215
              JWH-019, 1-hexyl-3-(1-naphthoyl)indole;
216
              JWH-073, 1-butyl-3-(1-naphthoyl)indole;
217
              JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)- Naphthalen-1-ylmethanone;
218
              JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.]
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219 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-220 ADB); 221 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB); 222 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-223 AMB); 224 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA); 225 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide 226 (ADB-FUBINACA); 227 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate 228 (MDMB-CHMICA); 229 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate Methyl 230 (MDMB-FUBINACA); 231 Tetrahydrocannabinols: 232 DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers. 233 DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers. 234 DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers. 235 Synthetic Phenethylamines 236 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-237 NBOMe); 238 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-239 NBOMe); 240 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-241 NBOMe); 242 Synthetic Opioids (icluding their isomers, esters, ethers, salts and salts of isomers, esters 243 and ethers): 244 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);

245	furanyl fentanyl;
246	3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
247	47700);
248	N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
249	phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);
250	N-[1-[2-hydroxy-2-(thiophen-2-yl)ethylpiperidin-4-yl]-N-phenylpropionamide, also known
251	as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-
252	hydroxythiofentanyl).
253	N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl)
254	N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl)
255	N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl)
256	2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide
257	(also known as U-48800)
258	Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as
259	U-49900)
260	Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also
261	known as U-51754)
262	Opioid Receptor Agonist
263	AH-7921 (3,4-dichloro-N- (1dimethylamino)cyclohexylmethyl]benzamide).
264	Naphthoylindoles or any compound containing a 3-(-1- Napthoyl) indole structure with
265	substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole
266	ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall
267	include the following:
268	JWH 015;
269	JWH 018;
270	JWH 019;

271	JWH 073;
272	JWH 081;
273	JWH 122;
274	JWH 200;
275	JWH 210;
276	JWH 398;
277	AM 2201;
278	WIN 55,212.
279	Naphylmethylindoles or any compound containing a 1hindol-3-yl-(1-naphthyl) methane
280	structure with a substitution at the nitrogen atom of the indole ring whether or not further
281	substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to
282	any extent. This shall include, but not be limited to, JWH 175 and JWH 184.
283	Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure with
284	substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the pyrrole
285	ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall
286	include, but not be limited to, JWH 147 and JWH 307.
287	Naphthylmethylindenes or any compound containing a Naphthylideneindene structure
288	with substitution at the 3- Position of the indene ring whether or not further substituted in the
289	indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This
290	shall include, but not be limited to, JWH 176.
291	Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure with
292	substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole
293	ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include
294	the following:
295	RCS-8, SR-18 OR BTM-8;
296	JWH 250;

297	JWH 203;
298	JWH 251;
299	JWH 302.
300	Cyclohexylphenols or any compound containing a 2-(3- hydroxycyclohexyl) phenol
301	structure with a substitution at the 5-position of the phenolic ring whether or not substituted in the
302	cyclohexyl ring to any extent. This shall include the following:
303	CP 47,497 and its homologues and analogs;
304	Cannabicyclohexanol;
305	CP 55,940.
306	Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with
307	substitution at the nitrogren atom of the indole ring whether or not further substituted in the indole
308	ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include
309	the following:
310	AM 694;
311	Pravadoline WIN 48,098;
312	RCS 4;
313	AM 679.
314	[2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-
315	napthalenymethanone. This shall include WIN 55,212-2.
316	Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol
317	structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051
318	and JWH 133.
319	Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with
320	substitution at the nitrogen atom of the indole ring whether or not further substituted in the
321	adamantoyl ring system to any extent. This shall include AM1248.

322	Tetramethylcyclopropylindoles or any compound containing A 3
323	tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring
324	whether or not further substituted in the indole ring to any extent and whether or not substituted
325	in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.
326	N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.
327	Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as
328	demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV
329	and V, not federal Food and Drug Administration approved drug or used within legitimate
330	approved medical research. Since nomenclature of these substances is not internationally
331	standardized, any immediate precursor or immediate derivative of these substances shall be
332	covered.
333	Tryptamines:
334	5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)
335	4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT)
336	4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)
337	4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)
338	4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT)
339	5-methoxy-α-methyltryptamine (5-MeO-AMT)
340	4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT)
341	4-hydroxy Diethyltryptamine (4-HO-DET)
342	5- methoxy- N,N- diallyltryptamine (5-MeO-DALT)
343	4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT)
344	4-hydroxy Diethyltryptamine (4-HO-DET)
345	FDU-PB-22 (1-Naphthyl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate);
346	FUB-PB-22 (Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate);
347	5-Fluoro-MN-24 (1-(5-Fluoropentyl)-N-(naphthalen-1-yl)-1H-indole-3-carboxamide);

348	MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
349	SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
350	SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
351	Methyl-Ethylaminopentiophenone;
352	FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
353	5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
354	5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
355	carboxamide);
356	MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
357	methylbutanoat);
358	MN-24 (N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide);
359	SDB-005 (Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate);
360	SDB-006 (1-Pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide);
361	Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
362	Methyl-Ethylaminopentiophenone;
363	FUB-AMB (Methyl(1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate);
364	5-Fluoro-SDB-005 Indole (Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate);
365	5F-AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-
366	3- carboxamide);
367	MMB-CHMICA (Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-
368	methylbutanoat);
369	Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
370	Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3-
371	a][1,4]benzodiazepine);
372	Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
373	benzodiazepin-2-one);

374	Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
375	a][1,4]diazepine);
376	Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
377	a][1,4]benzodiazepine);
378	Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
379	Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
380	a][1,4]benzodiazepine);
381	Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-
382	a][1,4]diazepine);
383	Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
384	one);
385	Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
386	Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
387	a][1,4]benzodiazepine).
388	(e) Depressants.
389	4-CN-CUMYL-BUTINACA (1-(4-Cyanobutyl)-N-(2-phenylpropan-2- yl)-1H-indazole-3-
390	carboxamide);
391	Alpha-Phenylacetoacetonitrile (3-Oxo-2-phenylbutanenitrile);
392	2-Fluoro Deschloroketamine (2-(2-Fluorophenyl)-2-(methylamino)-cyclohexanone,
393	monohydrochloride);
394	4-MEAP (2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one);
395	Mecloqualone;
396	Methaqualone.
397	Bromazolam (8-bromo-1-methyl-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
398	Clonazolam (6-(2-chlorophenyl)-1-methyl-8-nitro-4 H-[1,2,4]triazolo[4,3
399	a][1,4]benzodiazepine);

400	Cloniprazepam (5-(2-chlorophenyl)-1-(cyclopropylmethyl)-1,3-dihydro-7-nitro-2H-1,4-
401	benzodiazepin-2-one);
402	Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f] [1,2,4]triazolo[4,3-
403	a][1,4]diazepine);
404	Flualprazolam (8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
405	a][1,4]benzodiazepine);
406	Flubromazepam (7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one);
407	Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-
408	a][1,4]benzodiazepine);
409	Flunitrazolam (6-(2-fluorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-
410	a][1,4]diazepine);
411	Nifoxipam (5-(2-fluorophenyl)-1,3-dihydro-3-hydroxy-7-nitro-2H-1,4-benzodiazepin-2-
412	one);
413	Nitrazolam (1-methyl-8-nitro-6-phenyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine);
414	Pyrazolam (8-bromo-1-methyl-6-(2-pyridinyl)-4H-[1,2,4]triazolo[4,3-
415	a][1,4]benzodiazepine);
416	Declazepam (7-Chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-
417	one);
418	Deschloroetizolam (2-Ethyl-9-methyl-4-phenyl-6H-thieno[3,2- f][1,2,4]triazolo[4,3-
419	a][1,4]diazepine);
420	(f) Stimulants.
421	Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-
422	dihydro-5-phenyl-2-oxazolamine;
423	Cathinone; some trade or other names: 2-amino-1-phenyl-1- propanone, alpha-
424	aminopropiophenone, 2-aminopropiophenone and norephedrone;
425	Fenethylline:

426	Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
427	isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-
428	(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—-
429	methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or
430	mephedrone;3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone;
431	methylcathinone; AL-464; AL-422; AL- 463 and UR1432;
432	(+-) cis-4-methylaminorex; ((+-)cis-4,5-dihydro-4-methyl- 5-phenyl-2-oxazolamine);
433	N-ethylamphetamine;
434	N,N-dimethylamphetemine; also known as N,N-alpha- trimethyl-benzeneethanamine;
435	N,N-alpha-trimethylphenethylamine.
436	Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and
437	salts of isomers.
438	Substituted amphetamines:
439	2-Fluoroamphetamine
440	3-Fluoroamphetamine
441	4-Fluoroamphetamine
442	2-chloroamphetamine
443	3-chloroamphetamine
444	4-chloroamphetamine
445	2-Fluoromethamphetamine
446	3-Fluoromethamphetamine
447	4-Fluoromethamphetamine
448	4-chloromethamphetamine
449	Ethcathinone (2-(ethylamino)-1-phenyl-1-propanone, monohydrochloride);
450	Alpha-PHP (1-Phenyl-2-(pyrrolidin-1-yl)hexan-1-one);
451	MPHP (1-(4-Methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);

452	PV8 (1-Phenyl-2-(pyrrolidin-1-yl)heptan-1-one);
453	4-Chloro-Alpha-PVP (1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one);
454	N-Ethylhexedrone (2-(Ethylamino)-1-phenylhexan-1-one);
455	Methoxetamine (2-(Ethylamino)-2-(3-methoxyphenyl)-cyclohexanone);
456	3-Fluorophenmetrazine (2-(3-Fluorophenyl)-3-methylmorpholine):
457	(g) Temporary listing of substances subject to emergency scheduling. Any material,
458	compound, mixture or preparation which contains any quantity of the following substances:
459	N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts
460	and salts of isomers.
461	N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
462	isomers, salts and salts of isomers.
463	N-benzylpiperazine, also known as BZP.
464	Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);
465	4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-
466	butyramide);
467	Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);
468	Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-
469	acetamide);
470	3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-
471	phenylbutyramide);
472	4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
473	yl)butyramide);
474	Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);
475	Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
476	carboxamide);
477	Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).

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

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

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

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

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

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

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.

#### §60A-2-206. Schedule II.

- (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

9 indirectly by extraction from substances of vegetable origin, or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis: 10 11 Opium and opiate, and any salt, compound, derivative or preparation of opium or opiate 12 excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, 13 naloxone and naltrexone, and their respective salts, but including the following: 14 Raw opium; 15 Opium extracts; 16 Opium fluid; 17 Powdered opium; 18 Granulated opium; 19 Tincture of opium; 20 Codeine; 21 Dihydroetorphine; 22 Ethylmorphine; 23 Etorphine hydrochloride; 24 Hydrocodone; Hydromorphone; 25 26 Metopon; 27 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; Diphenoxylate; 51 52 Fentanyl; 53 Isomethadone; 54 Levo-alphacetylmethadol; 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	<u>Norfentanyl</u>
63	<u>Oliceridine</u>
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
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. —
80	Amphetamine, its salts, optical isomers and salts of its optical isomers;
81	Methamphetamine, its salts, isomers and salts of its isomers;
82	Methylphenidate;
83	Phenmetrazine and its salts; and
84	Lisdexamfetamine.
85	(e) Depressants. —
86	Amobarbital:

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4

87	Glutethimide;	
88	Pentobarbital;	
89	Phencyclidine;	
90	Secobarbital.	
91	(f) Hallucinogenic substances:	
92	Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] if in an FDA approved oral solution	
93	Nabilone: [Another name for nabilone: (+-)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8, 10,	
94	10a-hexahydro-1-hydroxy-6, 6-dimethyl-9H-dibenzo [b,d] pyran-9-one].	
95	(g) Immediate precursors. — Unless specifically excepted or unless listed in another	
96	schedule, any material, compound, mixture, or preparation which contains any quantity of the	
97	following substances:	
98	Immediate precursor to amphetamine and methamphetamine:	
99	Phenylacetone;	
100	Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl	
101	benzyl ketone;	
102	Immediate precursors to phencyclidine (PCP):	
103	1-phenylcyclohexylamine; and	
104	1-piperidinocyclohexanecarbonitrile (PCC).	
105	Immediate precursor to fentanyl:	
106	4-anilino-N-phenethyl-4-piperidine (ANPP).	
	§60A-2-210. Schedule IV.	
1	(a) Schedule IV shall consist of the drugs and other substances, by whatever official name,	
2	common or usual name, chemical name, or brand name designated, listed in this section. 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,

5	esters, ethers, saits and saits 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	Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-
14	propionoxybutane).
15	(c) Depressants.
16	Alprazolam;
17	Barbital;
18	Bromazepam;
19	Camazepam;
20	Carisoprodol;
21	Chloral betaine;
22	Chloral hydrate;
23	Chlordiazepoxide;
24	Clobazam;
25	Clonazepam;
26	Clorazepate;
27	Clotiazepam;
28	Cloxazolam;
29	Delorazepam;
30	Diazenam·

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	<u>Lemborexant</u> .
44	Loprazolam;
45	Lorazepam;
46	Lormetazepam;
47	Mebutamate;
48	Medazepam;
49	Meprobamate;
50	Methohexital;
51	Methylphenobarbital (mephobarbital);
52	Midazolam;
53	Nimetazepam;
54	Nitrazepam;
55	Nordiazepam;
56	Oxazepam;

57	Oxazolam;
58	Paraldehyde;
59	Petrichloral;
60	Phenobarbital;
61	Pinazepam;
62	Prazepam;
63	Quazepam;
64	Remimazolam.
65	Temazepam;
66	Tetrazepam;
67	Triazolam;
68	Zaleplon;
69	Zolpidem;
70	Zopiclone'
71	Suvorexant ([(7R)-4-(5-chloro-1,3-benzoxazol-2-yl)-7-methyl-1,4-diazepan-1-yl] [5-
72	methyl-2-(2H-1,2,3-triazol-2-yl)phenyl]methanone).
73	(d) Any material, compound, mixture or preparation which contains any quantity of
74	Fenfluramine and Dexfenfluramine.
75	(e) Stimulants.
76	Cathine ((+)-norpseudoephedrine);
77	Diethylpropion;
78	Fencamfamin;
79	Fenproporex;
80	Mazindol;
81	Mefenorex;
82	Modafinil;

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83	Pemoline (including organometallic complexes and chelates thereof);
84	Phentermine;
85	Pipradrol;
86	Serdexmethylphenidate.
87	Sibutramine;
88	SPA ((-)-1-dimethylamino-1,2-diphenylethane);
89	Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropyl
90	[(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid);
91	(f) Other substances. —
92	Pentazocine;
93	Butorphanol.
94	Tramadol (2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol);
95	Amyl nitrite, butyl nitrite, isobutyl nitrite and the other organic nitrites are controlled
96	substances and no product containing these compounds as a significant component shall be
97	possessed, bought or sold other than pursuant to a bona fide prescription or for industrial or
98	manufacturing purposes.
	§60A-2-212. Schedule V.
1	(a) Schedule V shall consist of the drugs and other substances, by whatever official name,
2	common or usual name, chemical name, or brand name designated, listed in this section. 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,

(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

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.

10	one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the
11	compound, mixture or preparation valuable medicinal qualities other than those possessed by the
12	narcotic drug alone.
13	Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;
14	Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams;
15	Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams;
16	Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine
17	sulfate per dosage unit;
18	Not more than 100 milligrams of opium per 100 milliliters or per 100 grams;
19	Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine
20	sulfate per dosage unit.
21	(c) Stimulants: —
22	Pyrovalerone.
23	(d) Any compound, mixture or preparation containing as its single active ingredient
24	ephedrine, pseudoephedrine or phenylpropanolamine, their salts or optical isomers, or salts of
25	optical isomers except products which are for pediatric use primarily intended for administration
26	to children under the age of 12: Provided, That neither the offenses set forth in section four
27	hundred one, article four of this chapter, nor the penalties therein, shall be applicable to ephedrine,
28	pseudoephedrine or phenylpropanolamine which shall be subject to the provisions of article ten
29	of this chapter.
30	(e) Depressants: —
31	Ezogabine [N-[2-amino-4-94-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester];
32	Lacosamide [(R)-2-acetoamido- N -benzyl-3-methoxy-propionamide];
33	Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid]; and
34	Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl] butanamide) (also referred to as
35	BRV; UCB-34714; Briviact).

- 36 (f) Other substances:
- 37 Gabapentin
- 38 Pregabalin