Amending Synthetic Drug Control Act

Summary

Amends ineffective law that simply address the problem of synthetic drugs by passing legislation to make specific chemical formulations of synthetic drugs illegal. These specific chemical formulations become dated quickly. This bill goes after broad base formulations that create or mimic the effect of cannabis or certain controlled substances and thereby helps law enforcement to stay ahead of criminal drug activity. Additionally it addresses weakness in the Analog statute that allows clever distributors to cloak products with an innocuous name in order to insulate themselves by marking packages: "Not For Human Consumption."

Model Legislation

Section 1. Short Title. Amending Synthetic Drug Control

Section 2. Purpose. Adds various synthetic drug compounds including various structural classes of those compounds to the list of Schedule I controlled substances. Deletes from the definition of controlled substance analog the requirement that a substance must be intended for human consumption to be considered an analog of a controlled substance with a chemical structure substantially similar to that of a controlled substance in Schedule I or II, or that was specifically designed to produce an effect substantially similar to that of a controlled substance in Schedule I or II.

Section 3. Definition.

(f) "Controlled Substance" means (i) a drug, substance, or immediate precursor, or synthetic drug in the Schedules of Article II of this Act or (ii) a drug or other substance, or immediate precursor, designated as a controlled substance by

the Department through administrative rule. The term does not include distilled spirits, wine, malt beverages, or tobacco, as terms are defined or used in... (Reference local state definitions of terms)

Section 4. Amending Synthetic Drug Control Act.

Sec. 204.

- (a) The controlled substances listed in this Section are included in Schedule I.
- (b) Unless specifically excepted or unless listed in another schedule, any of the following opiates, 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

45 designation:

- (1) Acetylmethadol;(1.1) Acetyl-alpha-methylfentanyl
- (N-[1-(1-methyl-2-phenethyl)-

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49
        4-piperidinyl]-N-phenylacetamide);
50
          (2) Allylprodine;
51
          (3) Alphacetylmethadol, except
52
        levo-alphacetylmethadol (also known as levo-alpha-
53
        acetylmethadol, levomethadyl acetate, or LAAM);
54
          (4) Alphameprodine;
55
          (5) Alphamethadol;
56
          (6) Alpha-methylfentanyl
57
        (N-(1-alpha-methyl-beta-phenyl) ethyl-4-piperidyl)
58
        propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-
59
        propanilido) piperidine;
60
          (6.1) Alpha-methylthiofentanyl
61
62
        (N-[1-methyl-2-(2-thienyl)ethyl-
63
        4-piperidinyl]-N-phenylpropanamide);
64
          (7) 1-methyl-4-phenyl-4-propionoxypiperidine (MPPP);
65
          (7.1) PEPAP
66
        (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
67
          (8) Benzethidine;
68
          (9) Betacetylmethadol;
69
          (9.1) Beta-hydroxyfentanyl
70
        (N-[1-(2-hydroxy-2-phenethyl)-
71
        4-piperidinyl]-N-phenylpropanamide);
72
          (10) Betameprodine;
73
          (11) Betamethadol;
74
          (12) Betaprodine;
75
          (13) Clonitazene;
76
          (14) Dextromoramide;
77
          (15) Diampromide;
78
          (16) Diethylthiambutene;
79
          (17) Difenoxin;
80
          (18) Dimenoxadol;
81
          (19) Dimepheptanol;
82
          (20) Dimethylthiambutene;
83
          (21) Dioxaphetylbutyrate;
84
          (22) Dipipanone;
85
          (23) Ethylmethylthiambutene;
86
          (24) Etonitazene;
87
          (25) Etoxeridine;
88
89
          (26) Furethidine;
90
          (27) Hydroxpethidine;
91
          (28) Ketobemidone;
92
          (29) Levomoramide;
93
          (30) Levophenacylmorphan;
94
          (31) 3-Methylfentanyl
95
        (N-[3-methyl-1-(2-phenylethyl)-
96
        4-piperidyl]-N-phenylpropanamide);
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97
            (31.1) 3-Methylthiofentanyl
 98
          (N-[(3-methyl-1-(2-thienyl)ethyl-
 99
          4-piperidinyl]-N-phenylpropanamide);
100
            (32) Morpheridine;
101
            (33) Noracymethadol;
102
            (34) Norlevorphanol;
103
            (35) Normethadone;
104
            (36) Norpipanone;
105
            (36.1) Para-fluorofentanyl
106
          (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-
107
          4-piperidinyl]propanamide);
108
            (37) Phenadoxone;
109
            (38) Phenampromide;
110
            (39) Phenomorphan;
111
            (40) Phenoperidine;
112
            (41) Piritramide;
113
            (42) Proheptazine;
            (43) Properidine;
114
115
116
            (44) Propiram;
117
            (45) Racemoramide;
118
            (45.1) Thiofentanyl
119
          (N-phenyl-N-[1-(2-thienyl)ethyl-
120
          4-piperidinyl]-propanamide);
121
            (46) Tilidine;
122
            (47) Trimeperidine;
123
            (48) Beta-hydroxy-3-methylfentanyl (other name:
124
          N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-
125
          N-phenylpropanamide).
126
          (c) Unless specifically excepted or unless listed in
127
        another schedule, any of the following opium derivatives, its
128
        salts, isomers and salts of isomers, whenever the existence of
129
        such salts, isomers and salts of isomers is possible within the
130
        specific chemical designation:
131
            (1) Acetorphine:
132
            (2) Acetyldihydrocodeine;
133
            (3) Benzylmorphine;
134
            (4) Codeine methylbromide;
135
            (5) Codeine-N-Oxide;
136
            (6) Cyprenorphine;
137
            (7) Desomorphine;
138
            (8) Diacetyldihydromorphine (Dihydroheroin);
139
            (9) Dihydromorphine;
140
            (10) Drotebanol;
141
            (11) Etorphine (except hydrochloride salt);
142
143
            (12) Heroin;
144
            (13) Hydromorphinol;
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145
            (14) Methyldesorphine:
146
            (15) Methyldihydromorphine;
147
            (16) Morphine methylbromide;
148
            (17) Morphine methylsulfonate;
149
            (18) Morphine-N-Oxide;
150
            (19) Myrophine;
151
            (20) Nicocodeine;
152
            (21) Nicomorphine;
153
            (22) Normorphine;
154
            (23) Pholcodine;
155
            (24) Thebacon.
156
          (d) Unless specifically excepted or unless listed in
157
       another schedule, any material, compound, mixture, or
158
       preparation which contains any quantity of the following
159
       hallucinogenic substances, or which contains any of its salts,
160
       isomers and salts of isomers, whenever the existence of such
161
       salts, isomers, and salts of isomers is possible within the
162
       specific chemical designation (for the purposes of this
163
       paragraph only, the term "isomer" includes the optical,
164
       position and geometric isomers):
165
            (1) 3,4-methylenedioxyamphetamine
166
          (alpha-methyl,3,4-methylenedioxyphenethylamine,
167
          methylenedioxyamphetamine, MDA);
168
            (1.1) Alpha-ethyltryptamine
169
170
          (some trade or other names: etryptamine;
171
          MONASE; alpha-ethyl-1H-indole-3-ethanamine;
172
          3-(2-aminobutyl)indole; a-ET; and AET);
173
            (2) 3,4-methylenedioxymethamphetamine (MDMA);
174
            (2.1) 3,4-methylenedioxy-N-ethylamphetamine
175
          (also known as: N-ethyl-alpha-methyl-
176
          3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE,
177
          and MDEA);
178
            (2.2) N-Benzylpiperazine (BZP);
179
            (2.2-1) Trifluoromethylphenylpiperazine (TFMPP);
180
            (3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);
181
            (4) 3,4,5-trimethoxyamphetamine (TMA);
182
            (5) (Blank);
183
            (6) Diethyltryptamine (DET);
184
            (7) Dimethyltryptamine (DMT);
185
            (7.1) 5-Methoxy-diallyltryptamine;
186
            (8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);
187
            (9) Ibogaine (some trade and other names:
188
          7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-
189
          6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b]
190
          indole; Tabernanthe iboga);
191
            (10) Lysergic acid diethylamide;
192
            (10.1) Salvinorin A;
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193
            (10.5) Salvia divinorum (meaning all parts of the plant
194
          presently classified botanically as Salvia divinorum,
195
          whether growing or not, the seeds thereof, any extract from
196
          any part of that plant, and every compound, manufacture,
197
          salts, isomers, and salts of isomers whenever the existence
198
          of such salts, isomers, and salts of isomers is possible
199
          within the specific chemical designation, derivative,
200
          mixture, or preparation of that plant, its seeds or
201
          extracts);
202
            (11) 3,4,5-trimethoxyphenethylamine (Mescaline);
203
            (12) Peyote (meaning all parts of the plant presently
204
          classified botanically as Lophophora williamsii Lemaire,
205
          whether growing or not, the seeds thereof, any extract from
206
          any part of that plant, and every compound, manufacture,
207
          salts, derivative, mixture, or preparation of that plant,
208
          its seeds or extracts);
209
            (13) N-ethyl-3-piperidyl benzilate (JB 318);
210
            (14) N-methyl-3-piperidyl benzilate;
211
            (14.1) N-hydroxy-3,4-methylenedioxyamphetamine
212
          (also known as N-hydroxy-alpha-methyl-
213
          3,4(methylenedioxy)phenethylamine and N-hydroxy MDA);
214
            (15) Parahexyl; some trade or other names:
215
          3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-
216
          dibenzo (b,d) pyran; Synhexyl;
217
            (16) Psilocybin;
218
            (17) Psilocyn;
219
            (18) Alpha-methyltryptamine (AMT);
220
            (19) 2,5-dimethoxyamphetamine
221
          (2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA);
222
223
            (20) 4-bromo-2,5-dimethoxyamphetamine
224
          (4-bromo-2,5-dimethoxy-alpha-methylphenethylamine;
225
          4-bromo-2,5-DMA);
226
            (20.1) 4-Bromo-2,5 dimethoxyphenethylamine.
227
          Some trade or other names: 2-(4-bromo-
228
          2,5-dimethoxyphenyl)-1-aminoethane;
229
          alpha-desmethyl DOB, 2CB, Nexus;
230
            (21) 4-methoxyamphetamine
231
          (4-methoxy-alpha-methylphenethylamine;
232
          paramethoxyamphetamine; PMA);
233
            (22) (Blank);
234
            (23) Ethylamine analog of phencyclidine.
235
          Some trade or other names:
236
          N-ethyl-1-phenylcyclohexylamine,
237
          (1-phenylcyclohexyl) ethylamine,
238
          N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE;
239
            (24) Pyrrolidine analog of phencyclidine. Some trade
240
          or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy,
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241
          PHP:
242
            (25) 5-methoxy-3,4-methylenedioxy-amphetamine;
243
            (26) 2,5-dimethoxy-4-ethylamphetamine
244
          (another name: DOET);
245
            (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine
246
          (another name: TCPy);
247
            (28) (Blank);
248
            (29) Thiophene analog of phencyclidine (some trade
249
250
          or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine;
251
          2-thienyl analog of phencyclidine; TPCP; TCP);
252
            (30) Bufotenine (some trade or other names:
253
          3-(Beta-Dimethylaminoethyl)-5-hydroxyindole;
254
          3-(2-dimethylaminoethyl)-5-indolol;
255
          5-hydroxy-N,N-dimethyltryptamine;
256
          N,N-dimethylserotonin; mappine);
257
            (31) 1-Pentyl-3-(1-naphthoyl)indole
258
          Some trade or other names: JWH-018;
259
            (32) 1-Butyl-3-(1-naphthoyl)indole
260
          Some trade or other names: JWH-073;
261
            (33) 1-[(5-fluoropentyl)-1H-indol-3-yl]-
262
          (2-iodophenyl)methanone
263
          Some trade or other names: AM-694;
264
            (34) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
265
          (2-methyloctan-2-yl)phenol
266
          Some trade or other names: CP 47,497
267
          and its C6, C8 and C9 homologs;
268
            (34.5) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
269
          (2-methyloctan-2-yl)phenol), where side chain n=5;
270
          and homologues where side chain n=4, 6, or 7; Some
271
          trade or other names: CP 47,497;
272
            (35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-
273
          (2-methyloctan-2-yl)-6a,7,
274
          10,10a-tetrahydrobenzo[c]chromen-1-ol
275
          Some trade or other names: HU-210;
276
277
            (35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-
278
          dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
279
          tetrahydrobenzo[c]chromen-1-ol, its isomers,
280
          salts, and salts of isomers; Some trade or other
281
          names: HU-210, Dexanabinol;
282
            (36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-
283
          6,6-dimethyl-3-(2-methyloctan-2-yl)-
284
          6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol
285
          Some trade or other names: HU-211;
286
            (37) (2-methyl-1-propyl-1H-indol-
287
          3-yl)-1-naphthalenyl-methanone
288
          Some trade or other names: JWH-015;
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289 (38) 4-methoxynaphthalen-1-yl-290 (1-pentylindol-3-yl)methanone 291 Some trade or other names: JWH-081; 292 (39) 1-Pentyl-3-(4-methyl-1-naphthoyl)indole 293 Some trade or other names: JWH-122; 294 (40) 2-(2-methylphenyl)-1-(1-pentyl-295 1H-indol-3-yl)-ethanone 296 Some trade or other names: JWH-251; 297 (41) 1-(2-cyclohexylethyl)-3-298 (2-methoxyphenylacetyl)indole 299 Some trade or other names: RCS-8, BTW-8 and SR-18; 300 (42) Any compound structurally derived from 301 3-(1-naphthoyl)indole or 1H-indol-3-vl-302 (1-naphthyl)methane by substitution at the 303 304 nitrogen atom of the indole ring by alkyl, haloalkyl, 305 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, 306 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, 307 or 2-(4-morpholinyl)ethyl whether or not further 308 substituted in the indole ring to any extent, whether 309 or not substituted in the naphthyl ring to any extent. 310 Examples of this structural class include, but are 311 not limited to, JWH-018, AM-2201, JWH-175, JWH-184, 312 and JWH-185; 313 (43) Any compound structurally derived from 314 3-(1-naphthoyl)pyrrole by substitution at the nitrogen 315 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl, 316 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl 317 aryl halide, 1-(N-methyl-2-piperidinyl)methyl, 318 or 2-(4-morpholinyl)ethyl, whether or not further 319 substituted in the pyrrole ring to any extent, whether 320 or not substituted in the naphthyl ring to any extent. 321 Examples of this structural class include, but are not 322 limited to, JWH-030, JWH-145, JWH-146, JWH-307, and 323 JWH-368; 324 (44) Any compound structurally derived from 325 1-(1-naphthylmethyl)indene by substitution 326 at the 3-position of the indene ring by alkyl, haloalkyl, 327 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl 328 halide, alkyl aryl halide, 1-(N-methyl-329 2-piperidinyl)methyl, or 2-(4-330 331 morpholinyl)ethyl whether or not further substituted in 332 the indene ring to any extent, whether or not substituted 333 in the naphthyl ring to any extent. Examples of 334 this structural class include, but are not 335 limited to, JWH-176; 336 (45) Any compound structurally derived from

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337
          3-phenylacetylindole by substitution at the
338
          nitrogen atom of the indole ring with alkyl, haloalkyl,
339
          alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl
340
          halide, alkyl aryl halide, 1-(N-methyl-2-
341
          piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
342
          whether or not further substituted in the indole ring
343
          to any extent, whether or not substituted in the phenyl
344
          ring to any extent. Examples of this structural
345
          class include, but are not limited to, JWH-167,
346
          JWH-250, JWH-251, and RCS-8;
347
            (46) Any compound structurally derived from
348
          2-(3-hydroxycyclohexyl)phenol by substitution
349
          at the 5-position of the phenolic ring by alkyl,
350
          haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
351
          aryl halide, alkyl aryl halide, 1-(N-methyl-2-
352
          piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
353
          whether or not substituted in the cyclohexyl ring to any
354
          extent. Examples of this structural class
355
          include, but are not limited to, CP 47,
356
          497 and its C8 homologue (cannabicyclohexanol);
357
358
            (46.1) Benzoylindoles: Any compound
359
          containing a 3-(benzoyl) indole structure with
360
          substitution at the nitrogen atom of the
361
          indole ring by an alkyl, haloalkyl, alkenyl,
362
          cycloalkylmethyl, cycloalkylethyl,
363
          1-(N-methyl-2-piperidinyl)methyl,
364
          or 2-(4-morpholinyl)ethyl group
365
          whether or not further substituted
366
          in the indole ring to any extent and
367
          whether or not substituted in the phenyl ring
368
          to any extent. Examples of this structural class
369
          include, but are not limited, to AM-630,
370
          AM-2233, AM-694, Pravadoline (WIN 48,098), and RCS-4;
371
            (47) 3,4-Methylenedioxymethcathinone
372
          Some trade or other names: Methylone;
373
            (48) 3,4-Methyenedioxypyrovalerone
374
          Some trade or other names: MDPV;
375
            (49) 4-Methylmethcathinone
376
          Some trade or other names: Mephedrone;
377
            (50) 4-methoxymethcathinone;
378
            (51) 4-Fluoromethcathinone;
379
            (52) 3-Fluoromethcathinone;
380
            (53) 2,5-Dimethoxy-4-(n)-propylthio-
381
          phenethylamine;
382
            (54) 5-Methoxy-N,N-diisopropyltryptamine;
383
            (55) Pentedrone:
384
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385 (56) 4-iodo-2,5-dimethoxy-N-((2-methoxy 386 phenyl)methyl)-benzeneethanamine 387 (trade or other name: 25I-NBOMe); 388 (57) 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl) 389 methyl]-benzeneethanamine (trade or other name: 390 25C-NBOMe); 391 (58) 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl) 392 methyl]-benzeneethanamine (trade or other name: 393 25B-NBOMe); . 394 (59) 3-cyclopropoylindole with 395 substitution at the nitrogen atom of the 396 indole ring by alkyl, haloalkyl, alkenyl, 397 cycloalkylmethyl, cycloalkylethyl, aryl 398 halide, alkyl aryl halide, 399 1-(N-methyl-2-piperidinyl)methyl, or 400 2-(4-morpholinyl)ethyl, whether or not 401 further substituted on the indole ring 402 to any extent, whether or not substituted 403 on the cyclopropyl ring to any extent: 404 including but not limited to XLR11, 405 UR144, FUB-144; 406 (60) 3-adamantoylindole with 407 substitution at the nitrogen atom of the 408 indole ring by alkyl, haloalkyl, alkenyl, 409 cycloalkylmethyl, cycloalkylethyl, 410 aryl halide, alkyl aryl halide, 411 412 1-(N-methyl-2-piperidinyl)methyl, or 413 2-(4-morpholinyl)ethyl, whether or not 414 further substituted on the indole ring to 415 any extent, whether or not substituted on 416 the adamantyl ring to any extent: including 417 but not limited to AB-001; 418 (61) N-(adamantyl)-indole-3-carboxamide 419 with substitution at the nitrogen atom of the 420 indole ring by alkyl, haloalkyl, alkenyl, 421 cycloalkylmethyl, cycloalkylethyl, aryl halide, 422 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, 423 or 2-(4-morpholinyl)ethyl, whether or not further 424 substituted on the indole ring to any extent, whether 425 or not substituted on the adamantyl ring to any 426 extent: including but not limited to 427 APICA/2NE-1, STS-135; 428 (62) N-(adamantyl)-indazole-3-carboxamide 429 with substitution at a nitrogen atom of the indazole 430 ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, 431 cycloalkylethyl, aryl halide, alkyl aryl halide, 432 1-(N-methyl-2-piperidinyl)methyl, or

2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the adamantyl ring to any extent: including but not limited to AKB48, 5F-AKB48;

- (63) 1H-indole-3-carboxylic acid 8-quinolinyl ester with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent, whether or not substituted on the quinoline ring to any extent: including but not limited to PB22, 5F-PB22, FUB-PB-22;
- (64) 3-(1-naphthoyl)indazole with substitution at the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the naphthyl ring to any extent: including but not limited to THJ-018, THJ-2201;
- (65) 2-(1-naphthoyl)benzimidazole with substitution at the nitrogen atom of the benzimidazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or

- 2-(4-morpholinyl)ethyl, whether or not further substituted on the benzimidazole ring to any extent, whether or not substituted on the naphthyl ring to any extent: including, but not limited to FUBIMINA;
- (66) N-(1-amino-3-methyl-1-oxobutan-2-yl)
 -1H-indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including but not limited to AB-PINACA, AB-FUBINACA, AB-CHMINACA;
- (67) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indazole-3-carboxamide with substitution

on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including but not limited to ADB-PINACA, ADB-FUBINACA;

(68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl

aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent: including but not limited to ADBICA, 5F-ADBICA;

(69) N-(1-amino-3-methyl-1-oxobutan-2-yl)1H-indole-3-carboxamide with substitution on the
nitrogen atom of the indole ring by alkyl, haloalkyl,
alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl
halide, alkyl aryl halide, 1-(N-methyl-2piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
whether or not further substituted on the indole
ring to any extent: including but not limited
to ABICA, 5F-ABICA;

(70) Methyl 2-(1H-indazole-3-carboxamido)-3-methylbutanoate with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including but not limited to AMB, 5F-AMB.

(e) Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous

system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) mecloqualone;
- (2) methagualone; and
- (3) gamma hydroxybutyric acid.
- (f) Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following

529 substances having a stimulant effect on the central nervous 530 system, including its salts, isomers, and salts of isomers: 531 (1) Fenethylline; 532 (2) N-ethylamphetamine; 533 (3) Aminorex (some other names: 534 2-amino-5-phenyl-2-oxazoline; aminoxaphen; 535 4-5-dihydro-5-phenyl-2-oxazolamine) and its 536 salts, optical isomers, and salts of optical isomers; 537 (4) Methcathinone (some other names: 538 2-methylamino-1-phenylpropan-1-one; 539 Ephedrone; 2-(methylamino)-propiophenone; 540 alpha-(methylamino)propiophenone; N-methylcathinone; 541 methycathinone; Monomethylpropion; UR 1431) and its 542 salts, optical isomers, and salts of optical isomers; 543 (5) Cathinone (some trade or other names: 544 2-aminopropiophenone; alpha-aminopropiophenone; 545 2-amino-1-phenyl-propanone; norephedrone); 546 547 (6) N,N-dimethylamphetamine (also known as: 548 N,N-alpha-trimethyl-benzeneethanamine; 549 N,N-alpha-trimethylphenethylamine); 550 (7) (+ or -) cis-4-methylaminorex ((+ or -) cis-551 4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine); 552 (8) 3,4-Methylenedioxypyrovalerone (MDPV). 553 (g) Temporary listing of substances subject to emergency 554 scheduling. Any material, compound, mixture, or preparation 555 that contains any quantity of the following substances: 556 (1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide 557 (benzylfentanyl), its optical isomers, isomers, salts, 558 and salts of isomers; 559 (2) N-[1(2-thienyl) 560 methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), 561 its optical isomers, salts, and salts of isomers. 562 (h) Synthetic cathinones. Unless specifically excepted, 563 any chemical compound not including bupropion, structurally 564 derived from 2-aminopropan-1-one by substitution at the 565 1-position with either phenyl, naphthyl, or thiophene ring 566 systems, whether or not the compound is further modified in one 567 or more of the following ways: 568 (1) by substitution in the ring system to 569 any extent with alkyl, alkylenedioxy, alkoxy, 570 haloalkyl, hydroxyl, or halide substituents, whether 571 or not further substituted in the ring system 572 by one or more other univalent substituents. 573 574 Examples of this class include, but are not 575 limited to, 3,4-Methylenedioxycathinone 576 (bk-MDA);

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577
            (2) by substitution at the 3-position
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          with an acyclic alkyl substituent. Examples of
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          this class include, but are not limited to,
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          2-methylamino-1-phenylbutan-1-one
581
          (buphedrone); or
582
            (3) by substitution at the 2-amino nitrogen
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          atom with alkyl, dialkyl, benzyl, or methoxybenzyl
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          groups, or by inclusion of the 2-amino nitrogen atom
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          in a cyclic structure. Examples of this class include,
586
          but are not limited to, Dimethylcathinone, Ethcathinone,
587
          and a-Pyrrolidinopropiophenone (a-PPP).
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        (Source: P.A. 97-192, eff. 7-22-11; 97-193, eff. 1-1-12;
589
        97-194, eff. 7-22-11; 97-334, eff. 1-1-12; 97-813, eff.
590
        7-13-12; 97-872, eff. 7-31-12; 98-987, eff. 1-1-15.)
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592
        Sec (Regarding Manufacture or Delivery of Controlled Synthetic drugs). 401.
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        Except as authorized by this Act, it is unlawful
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        for any person knowingly to manufacture or deliver, or possess
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        with intent to manufacture or deliver, a controlled substance
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        other than methamphetamine, a counterfeit substance, or a
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        controlled substance analog. A violation of this Act with
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        respect to each of the controlled substances listed herein
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        constitutes a single and separate violation of this Act. For
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        purposes of this Section, "controlled substance analog" or
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        "analog" means a substance which is intended for human
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        consumption, other than a controlled substance, that has a
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        chemical structure substantially similar to that of a
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        controlled substance in Schedule I or II, or that was
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        specifically designed to produce an effect substantially
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        similar to that of a controlled substance in Schedule I or II.
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        Examples of chemical classes in which controlled substance
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        analogs are found include, but are not limited to, the
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        following: phenethylamines, N-substituted piperidines,
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        morphinans, ecgonines, quinazolinones, substituted indoles,
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        and arylcycloalkylamines. For purposes of this Act, a
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        controlled substance analog shall be treated in the same manner
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        as the controlled substance to which it is substantially
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        similar.
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616
        Sec. 402 (Regarding Possession of Controlled Synthetic Drugs).
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        Except as otherwise authorized by this Act, it is
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        unlawful for any person knowingly to possess a controlled or
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        counterfeit substance or controlled substance analog. A
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        violation of this Act with respect to each of the controlled
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        substances listed herein constitutes a single and separate
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        violation of this Act. For purposes of this Section,
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        "controlled substance analog" or "analog" means a substance
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        which is intended for human consumption, other than a
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625 controlled substance, that has a chemical structure 626 substantially similar to that of a controlled substance in 627 Schedule I or II, or that was specifically designed to produce 628 629 an effect substantially similar to that of a controlled 630 substance in Schedule I or II. Examples of chemical classes in 631 which controlled substance analogs are found include, but are 632 not limited to, the following: phenethylamines, N-substituted 633 piperidines, morphinans, ecgonines, quinazolinones, 634 substituted indoles, and arylcycloalkylamines. For purposes of 635 this Act, a controlled substance analog shall be treated in the 636 same manner as the controlled substance to which it is 637 substantially similar. 638 639 Section 5. Severability clause. 640 641 Section 6. Repealer clause. 642 643 **Section 7. {Insert Effective Date}**