

By the Committee on Fiscal Policy; and Senator Yarborough

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1 A bill to be entitled
2 An act relating to intoxicating substances; providing
3 a short title; creating s. 569.216, F.S.; prohibiting
4 tobacco or nicotine dealers, or their agents or
5 employees, from possessing, selling, possessing with
6 intent to sell, delivering, or giving, directly or
7 indirectly, nitrous oxide on or from the dealer's
8 licensed premises; providing criminal penalties;
9 providing applicability; providing an exception;
10 requiring the Department of Business and Professional
11 Regulation to adopt rules; amending s. 893.03, F.S.;
12 excepting from the list of Schedule I controlled
13 substances certain xylazine animal drug products
14 approved by the United States Food and Drug
15 Administration and used for certain purposes; amending
16 s. 893.13, F.S.; providing criminal penalties and
17 requiring a mandatory minimum term of imprisonment if
18 a person sells, manufactures, or delivers or possesses
19 with intent to sell, manufacture, or deliver xylazine;
20 amending s. 893.135, F.S.; creating the offense of
21 trafficking in xylazine; providing criminal penalties
22 and requiring a mandatory minimum term of imprisonment
23 and fines based on the quantity of the controlled
24 substance involved in the offense; providing effective
25 dates.

26
27 Be It Enacted by the Legislature of the State of Florida:

28
29 Section 1. Section 2 of this act may be cited as "Meg's

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30 Law."

31 Section 2. Section 569.216, Florida Statutes, is created to
32 read:

33 569.216 Prohibition on possessing, selling, furnishing, or
34 giving nitrous oxide; exceptions; penalties.—

35 (1) It is unlawful for any dealer who is licensed or
36 permitted under this chapter, or a dealer's agent or employee,
37 to possess, sell, possess with intent to sell, deliver, or give,
38 directly or indirectly, nitrous oxide on or from the dealer's
39 licensed premises. A dealer or a dealer's agent or employee who
40 violates this subsection commits a felony of the third degree,
41 punishable as provided in s. 775.082, s. 775.083, or s. 775.084.

42 (2) This section does not apply to a grocery store or
43 supermarket, as licensed or permitted by the Department of
44 Agriculture and Consumer Services, but does apply to a
45 convenience business, as defined by s. 812.171.

46 (3) This section does not prohibit the purchase or sale of
47 a finished food product in which nitrous oxide is used solely as
48 a propellant.

49 (4) The Department of Business and Professional Regulation
50 shall adopt rules regarding the sale and purchase of nitrous
51 oxide to prevent the use of nitrous oxide for inducing a
52 condition of intoxication. Such rules may address products
53 containing nitrous oxide and finished food products in which
54 nitrous oxide is used solely as a propellant.

55 Section 3. Effective July 1, 2026, paragraph (c) of
56 subsection (1) of section 893.03, Florida Statutes, is amended
57 to read:

58 893.03 Standards and schedules.—The substances enumerated

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59 in this section are controlled by this chapter. The controlled
60 substances listed or to be listed in Schedules I, II, III, IV,
61 and V are included by whatever official, common, usual,
62 chemical, trade name, or class designated. The provisions of
63 this section shall not be construed to include within any of the
64 schedules contained in this section any excluded drugs listed
65 within the purview of 21 C.F.R. s. 1308.22, styled "Excluded
66 Substances"; 21 C.F.R. s. 1308.24, styled "Exempt Chemical
67 Preparations"; 21 C.F.R. s. 1308.32, styled "Exempted
68 Prescription Products"; or 21 C.F.R. s. 1308.34, styled "Exempt
69 Anabolic Steroid Products."

70 (1) SCHEDULE I.—A substance in Schedule I has a high
71 potential for abuse and has no currently accepted medical use in
72 treatment in the United States and in its use under medical
73 supervision does not meet accepted safety standards. The
74 following substances are controlled in Schedule I:

75 (c) Unless specifically excepted or unless listed in
76 another schedule, any material, compound, mixture, or
77 preparation that contains any quantity of the following
78 hallucinogenic substances or that contains any of their salts,
79 isomers, including optical, positional, or geometric isomers,
80 homologues, nitrogen-heterocyclic analogs, esters, ethers, and
81 salts of isomers, homologues, nitrogen-heterocyclic analogs,
82 esters, or ethers, if the existence of such salts, isomers, and
83 salts of isomers is possible within the specific chemical
84 designation or class description:

- 85 1. Alpha-Ethyltryptamine.
- 86 2. 4-Methylaminorex (2-Amino-4-methyl-5-phenyl-2-
87 oxazoline).

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- 88 3. Aminorex (2-Amino-5-phenyl-2-oxazoline).
89 4. DOB (4-Bromo-2,5-dimethoxyamphetamine).
90 5. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
91 6. Bufotenine.
92 7. Cannabis.
93 8. Cathinone.
94 9. DET (Diethyltryptamine).
95 10. 2,5-Dimethoxyamphetamine.
96 11. DOET (4-Ethyl-2,5-Dimethoxyamphetamine).
97 12. DMT (Dimethyltryptamine).
98 13. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine
99 analog of phencyclidine).
100 14. JB-318 (N-Ethyl-3-piperidyl benzilate).
101 15. N-Ethylamphetamine.
102 16. Fenethylamine.
103 17. 3,4-Methylenedioxy-N-hydroxyamphetamine.
104 18. Ibogaine.
105 19. LSD (Lysergic acid diethylamide).
106 20. Mescaline.
107 21. Methcathinone.
108 22. 5-Methoxy-3,4-methylenedioxyamphetamine.
109 23. PMA (4-Methoxyamphetamine).
110 24. PMMA (4-Methoxymethamphetamine).
111 25. DOM (4-Methyl-2,5-dimethoxyamphetamine).
112 26. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
113 27. MDA (3,4-Methylenedioxyamphetamine).
114 28. JB-336 (N-Methyl-3-piperidyl benzilate).
115 29. N,N-Dimethylamphetamine.
116 30. Parahexyl.

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- 117 31. Peyote.
- 118 32. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine
119 analog of phencyclidine).
- 120 33. Psilocybin.
- 121 34. Psilocyn.
- 122 35. *Salvia divinorum*, except for any drug product approved
123 by the United States Food and Drug Administration which contains
124 *Salvia divinorum* or its isomers, esters, ethers, salts, and
125 salts of isomers, esters, and ethers, if the existence of such
126 isomers, esters, ethers, and salts is possible within the
127 specific chemical designation.
- 128 36. Salvinorin A, except for any drug product approved by
129 the United States Food and Drug Administration which contains
130 Salvinorin A or its isomers, esters, ethers, salts, and salts of
131 isomers, esters, and ethers, if the existence of such isomers,
132 esters, ethers, and salts is possible within the specific
133 chemical designation.
- 134 37. Xylazine, except for a xylazine animal drug product
135 approved by the United States Food and Drug Administration and
136 the use of which conforms to the approved application or is
137 authorized under 21 U.S.C. s. 360b(a)(4). The manufacture,
138 importation, distribution, prescribing, or sale of xylazine for
139 human use is not subject to this exception.
- 140 38. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine)
141 (Thiophene analog of phencyclidine).
- 142 39. 3,4,5-Trimethoxyamphetamine.
- 143 40. Methydone (3,4-Methylenedioxymethcathinone).
- 144 41. MDPV (3,4-Methylenedioxypyrovalerone).
- 145 42. Methyldmethcathinone.

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- 146 43. Methoxymethcathinone.
- 147 44. Fluoromethcathinone.
- 148 45. Methylethcathinone.
- 149 46. CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-
150 yl)phenol) and its dimethyloctyl (C8) homologue.
- 151 47. HU-210 [(6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-
152 methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol].
- 153 48. JWH-018 (1-Pentyl-3-(1-naphthoyl)indole).
- 154 49. JWH-073 (1-Butyl-3-(1-naphthoyl)indole).
- 155 50. JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-
156 naphthoyl)indole).
- 157 51. BZP (Benzylpiperazine).
- 158 52. Fluorophenylpiperazine.
- 159 53. Methylphenylpiperazine.
- 160 54. Chlorophenylpiperazine.
- 161 55. Methoxyphenylpiperazine.
- 162 56. DBZP (1,4-Dibenzylpiperazine).
- 163 57. TFMPP (Trifluoromethylphenylpiperazine).
- 164 58. MBDB (Methylbenzodioxolylbutanamine) or (3,4-
165 Methylenedioxy-N-methylbutanamine).
- 166 59. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
- 167 60. 5-Hydroxy-N-methyltryptamine.
- 168 61. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).
- 169 62. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
- 170 63. Methyltryptamine.
- 171 64. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
- 172 65. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
- 173 66. Tyramine (4-Hydroxyphenethylamine).
- 174 67. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).

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- 175 68. DiPT (N,N-Diisopropyltryptamine).
176 69. DPT (N,N-Dipropyltryptamine).
177 70. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
178 71. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
179 72. DOI (4-Iodo-2,5-dimethoxyamphetamine).
180 73. DOC (4-Chloro-2,5-dimethoxyamphetamine).
181 74. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
182 75. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
183 76. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
184 77. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
185 78. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
186 79. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
187 80. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
188 81. Butylone (3,4-Methylenedioxy-alpha-
189 methylaminobutyrophenone).
190 82. Ethcathinone.
191 83. Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
192 84. Naphyrone (Naphthylpyrovalerone).
193 85. Dimethylone (3,4-Methylenedioxy-N,N-dimethylcathinone).
194 86. 3,4-Methylenedioxy-N,N-diethylcathinone.
195 87. 3,4-Methylenedioxy-propiofenone.
196 88. 3,4-Methylenedioxy-alpha-bromopropiofenone.
197 89. 3,4-Methylenedioxy-propiofenone-2-oxime.
198 90. 3,4-Methylenedioxy-N-acetylcathinone.
199 91. 3,4-Methylenedioxy-N-acetylmethcathinone.
200 92. 3,4-Methylenedioxy-N-acetylethcathinone.
201 93. Bromomethcathinone.
202 94. Buphedrone (alpha-Methylamino-butyrophenone).
203 95. Eutylone (3,4-Methylenedioxy-alpha-

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204 ethylaminobutyrophenone).

205 96. Dimethylcathinone.

206 97. Dimethylmethcathinone.

207 98. Pentylone (3,4-Methylenedioxy-alpha-

208 methylaminovalerophenone).

209 99. MDPMP (3,4-Methylenedioxy-alpha-

210 pyrrolidinopropiophenone).

211 100. MDPPB (3,4-Methylenedioxy-alpha-

212 pyrrolidinobutyrophenone).

213 101. MOPPP (Methoxy-alpha-pyrrolidinopropiophenone).

214 102. MPHP (Methyl-alpha-pyrrolidinohexanophenone).

215 103. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP

216 (Benocyclidine).

217 104. F-MABP (Fluoromethylaminobutyrophenone).

218 105. MeO-PBP (Methoxypyrrolidinobutyrophenone).

219 106. Et-PBP (Ethylpyrrolidinobutyrophenone).

220 107. 3-Me-4-MeO-MCAT (3-Methyl-4-Methoxymethcathinone).

221 108. Me-EABP (Methylethylaminobutyrophenone).

222 109. Etizolam.

223 110. PPP (Pyrrolidinopropiophenone).

224 111. PBP (Pyrrolidinobutyrophenone).

225 112. PVP (Pyrrolidinovalerophenone) or

226 (Pyrrolidinopentiophenone).

227 113. MPPP (Methyl-alpha-pyrrolidinopropiophenone).

228 114. JWH-007 (1-Pentyl-2-methyl-3-(1-naphthoyl)indole).

229 115. JWH-015 (1-Propyl-2-methyl-3-(1-naphthoyl)indole).

230 116. JWH-019 (1-Hexyl-3-(1-naphthoyl)indole).

231 117. JWH-020 (1-Heptyl-3-(1-naphthoyl)indole).

232 118. JWH-072 (1-Propyl-3-(1-naphthoyl)indole).

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- 233 119. JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl)indole).
- 234 120. JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
- 235 121. JWH-133 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-methylpentan-
- 236 2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).
- 237 122. JWH-175 (1-Pentyl-3-(1-naphthylmethyl)indole).
- 238 123. JWH-201 (1-Pentyl-3-(4-methoxyphenylacetyl)indole).
- 239 124. JWH-203 (1-Pentyl-3-(2-chlorophenylacetyl)indole).
- 240 125. JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).
- 241 126. JWH-250 (1-Pentyl-3-(2-methoxyphenylacetyl)indole).
- 242 127. JWH-251 (1-Pentyl-3-(2-methylphenylacetyl)indole).
- 243 128. JWH-302 (1-Pentyl-3-(3-methoxyphenylacetyl)indole).
- 244 129. JWH-398 (1-Pentyl-3-(4-chloro-1-naphthoyl)indole).
- 245 130. HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-
- 246 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-
- 247 ol).
- 248 131. HU-308 ([(1R,2R,5R) -2-[2,6-Dimethoxy-4-(2-methyloctan-
- 249 2-yl)phenyl]-7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl]
- 250 methanol).
- 251 132. HU-331 (3-Hydroxy-2-[(1R,6R)-3-methyl-6-(1-
- 252 methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-2,5-cyclohexadiene-
- 253 1,4-dione).
- 254 133. CB-13 (4-Pentyloxy-1-(1-naphthoyl)naphthalene).
- 255 134. CB-25 (N-Cyclopropyl-11-(3-hydroxy-5-pentylphenoxy)-
- 256 undecanamide).
- 257 135. CB-52 (N-Cyclopropyl-11-(2-hexyl-5-hydroxyphenoxy)-
- 258 undecanamide).
- 259 136. CP 55,940 (2-[3-Hydroxy-6-propanol-cyclohexyl]-5-(2-
- 260 methyloctan-2-yl)phenol).
- 261 137. AM-694 (1-(5-Fluoropentyl)-3-(2-iodobenzoyl)indole).

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- 262 138. AM-2201 (1-(5-Fluoropentyl)-3-(1-naphthoyl)indole).
- 263 139. RCS-4 (1-Pentyl-3-(4-methoxybenzoyl)indole).
- 264 140. RCS-8 (1-(2-Cyclohexylethyl)-3-(2-
- 265 methoxyphenylacetyl)indole).
- 266 141. WIN55,212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-
- 267 morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-
- 268 naphthalenylmethanone).
- 269 142. WIN55,212-3 ([(3S)-2,3-Dihydro-5-methyl-3-(4-
- 270 morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-
- 271 naphthalenylmethanone).
- 272 143. Pentedrone (alpha-Methylaminovalerophenone).
- 273 144. Fluoroamphetamine.
- 274 145. Fluoromethamphetamine.
- 275 146. Methoxetamine.
- 276 147. Methiopropamine.
- 277 148. Methylbuphedrone (Methyl-alpha-
- 278 methylaminobutyrophenone).
- 279 149. APB ((2-Aminopropyl)benzofuran).
- 280 150. APDB ((2-Aminopropyl)-2,3-dihydrobenzofuran).
- 281 151. UR-144 (1-Pentyl-3-(2,2,3,3-
- 282 tetramethylcyclopropanoyl)indole).
- 283 152. XLR11 (1-(5-Fluoropentyl)-3-(2,2,3,3-
- 284 tetramethylcyclopropanoyl)indole).
- 285 153. Chloro UR-144 (1-(Chloropentyl)-3-(2,2,3,3-
- 286 tetramethylcyclopropanoyl)indole).
- 287 154. AKB48 (N-Adamant-1-yl 1-pentylindazole-3-carboxamide).
- 288 155. AM-2233(1-[(N-Methyl-2-piperidinyl)methyl]-3-(2-
- 289 iodobenzoyl)indole).
- 290 156. STS-135 (N-Adamant-1-yl 1-(5-fluoropentyl)indole-3-

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291 carboxamide).

292 157. URB-597 ((3'-(Aminocarbonyl)[1,1'-biphenyl]-3-yl)-

293 cyclohexylcarbamate).

294 158. URB-602 ([1,1'-Biphenyl]-3-yl-carbamic acid,

295 cyclohexyl ester).

296 159. URB-754 (6-Methyl-2-[(4-methylphenyl)amino]-1-

297 benzoxazin-4-one).

298 160. 2C-D (4-Methyl-2,5-dimethoxyphenethylamine).

299 161. 2C-H (2,5-Dimethoxyphenethylamine).

300 162. 2C-N (4-Nitro-2,5-dimethoxyphenethylamine).

301 163. 2C-P (4-(n)-Propyl-2,5-dimethoxyphenethylamine).

302 164. 25I-NBOMe (4-Iodo-2,5-dimethoxy-[N-(2-

303 methoxybenzyl)]phenethylamine).

304 165. MDMA (3,4-Methylenedioxymethamphetamine).

305 166. PB-22 (8-Quinolinyll 1-pentylindole-3-carboxylate).

306 167. Fluoro PB-22 (8-Quinolinyll 1-(fluoropentyl)indole-3-

307 carboxylate).

308 168. BB-22 (8-Quinolinyll 1-(cyclohexylmethyl)indole-3-

309 carboxylate).

310 169. Fluoro AKB48 (N-Adamant-1-yl 1-(fluoropentyl)indazole-

311 3-carboxamide).

312 170. AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-

313 pentylindazole-3-carboxamide).

314 171. AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-

315 (4-fluorobenzyl)indazole-3-carboxamide).

316 172. ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-

317 1-pentylindazole-3-carboxamide).

318 173. Fluoro ADBICA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-

319 yl)-1-(fluoropentyl)indole-3-carboxamide).

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- 320 174. 25B-NBOMe (4-Bromo-2,5-dimethoxy-[N-(2-
321 methoxybenzyl)]phenethylamine).
- 322 175. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-
323 methoxybenzyl)]phenethylamine).
- 324 176. AB-CHMINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-
325 (cyclohexylmethyl)indazole-3-carboxamide).
- 326 177. FUB-PB-22 (8-Quinolinyll 1-(4-fluorobenzyl)indole-3-
327 carboxylate).
- 328 178. Fluoro-NNEI (N-Naphthalen-1-yl 1-(fluoropentyl)indole-
329 3-carboxamide).
- 330 179. Fluoro-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-
331 (fluoropentyl)indazole-3-carboxamide).
- 332 180. THJ-2201 (1-(5-Fluoropentyl)-3-(1-naphthoyl)indazole).
- 333 181. AM-855 ((4aR,12bR)-8-Hexyl-2,5,5-trimethyl-
334 1,4,4a,8,9,10,11,12b-octahydronaphtho[3,2-c]isochromen-12-ol).
- 335 182. AM-905 ((6aR,9R,10aR)-3-[(E)-Hept-1-enyl]-9-
336 (hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-
337 hexahydrobenzo[c]chromen-1-ol).
- 338 183. AM-906 ((6aR,9R,10aR)-3-[(Z)-Hept-1-enyl]-9-
339 (hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-
340 hexahydrobenzo[c]chromen-1-ol).
- 341 184. AM-2389 ((6aR,9R,10aR)-3-(1-Hexyl-cyclobut-1-yl)-
342 6a,7,8,9,10,10a-hexahydro-6,6-dimethyl-6H-dibenzo[b,d]pyran-1,9
343 diol).
- 344 185. HU-243 ((6aR,8S,9S,10aR)-9-(Hydroxymethyl)-6,6-
345 dimethyl-3-(2-methyloctan-2-yl)-8,9-ditritio-7,8,10,10a-
346 tetrahydro-6aH-benzo[c]chromen-1-ol).
- 347 186. HU-336 ((6aR,10aR)-6,6,9-Trimethyl-3-pentyl-
348 6a,7,10,10a-tetrahydro-1H-benzo[c]chromene-1,4(6H)-dione).

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- 349 187. MAPB ((2-Methylaminopropyl)benzofuran).
350 188. 5-IT (2-(1H-Indol-5-yl)-1-methyl-ethylamine).
351 189. 6-IT (2-(1H-Indol-6-yl)-1-methyl-ethylamine).
352 190. Synthetic Cannabinoids.—Unless specifically excepted
353 or unless listed in another schedule or contained within a
354 pharmaceutical product approved by the United States Food and
355 Drug Administration, any material, compound, mixture, or
356 preparation that contains any quantity of a synthetic
357 cannabinoid found to be in any of the following chemical class
358 descriptions, or homologues, nitrogen-heterocyclic analogs,
359 isomers (including optical, positional, or geometric), esters,
360 ethers, salts, and salts of homologues, nitrogen-heterocyclic
361 analogs, isomers, esters, or ethers, whenever the existence of
362 such homologues, nitrogen-heterocyclic analogs, isomers, esters,
363 ethers, salts, and salts of isomers, esters, or ethers is
364 possible within the specific chemical class or designation.
365 Since nomenclature of these synthetically produced cannabinoids
366 is not internationally standardized and may continually evolve,
367 these structures or the compounds of these structures shall be
368 included under this subparagraph, regardless of their specific
369 numerical designation of atomic positions covered, if it can be
370 determined through a recognized method of scientific testing or
371 analysis that the substance contains properties that fit within
372 one or more of the following categories:
- 373 a. Tetrahydrocannabinols.—Any tetrahydrocannabinols
374 naturally contained in a plant of the genus *Cannabis*, the
375 synthetic equivalents of the substances contained in the plant
376 or in the resinous extracts of the genus *Cannabis*, or synthetic
377 substances, derivatives, and their isomers with similar chemical

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378 structure and pharmacological activity, including, but not
379 limited to, Delta 9 tetrahydrocannabinols and their optical
380 isomers, Delta 8 tetrahydrocannabinols and their optical
381 isomers, Delta 6a,10a tetrahydrocannabinols and their optical
382 isomers, or any compound containing a tetrahydrobenzo[c]chromene
383 structure with substitution at either or both the 3-position or
384 9-position, with or without substitution at the 1-position with
385 hydroxyl or alkoxy groups, including, but not limited to:

386 (I) Tetrahydrocannabinol.

387 (II) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-
388 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-
389 ol).

390 (III) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-
391 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-
392 ol).

393 (IV) JWH-051 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-
394 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

395 (V) JWH-133 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-methylpentan-
396 2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

397 (VI) JWH-057 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-methyloctan-
398 2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

399 (VII) JWH-359 ((6aR,10aR)-1-Methoxy-6,6,9-trimethyl-3-(2,3-
400 dimethylpentan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

401 (VIII) AM-087 ((6aR,10aR)-3-(2-Methyl-6-bromohex-2-yl)-
402 6,6,9-trimethyl-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol).

403 (IX) AM-411 ((6aR,10aR)-3-(1-Adamantyl)-6,6,9-trimethyl-
404 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol).

405 (X) Parahexyl.

406 b. Naphthoylindoles, Naphthoylindazoles,

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407 Naphthoylcarbazoles, Naphthylmethylindoles,
408 Naphthylmethylindazoles, and Naphthylmethylcarbazoles.—Any
409 compound containing a naphthoylindole, naphthoylindazole,
410 naphthoylcarbazole, naphthylmethylindole,
411 naphthylmethylindazole, or naphthylmethylcarbazole structure,
412 with or without substitution on the indole, indazole, or
413 carbazole ring to any extent, whether or not substituted on the
414 naphthyl ring to any extent, including, but not limited to:

- 415 (I) JWH-007 (1-Pentyl-2-methyl-3-(1-naphthoyl)indole).
- 416 (II) JWH-011 (1-(1-Methylhexyl)-2-methyl-3-(1-
417 naphthoyl)indole).
- 418 (III) JWH-015 (1-Propyl-2-methyl-3-(1-naphthoyl)indole).
- 419 (IV) JWH-016 (1-Butyl-2-methyl-3-(1-naphthoyl)indole).
- 420 (V) JWH-018 (1-Pentyl-3-(1-naphthoyl)indole).
- 421 (VI) JWH-019 (1-Hexyl-3-(1-naphthoyl)indole).
- 422 (VII) JWH-020 (1-Heptyl-3-(1-naphthoyl)indole).
- 423 (VIII) JWH-022 (1-(4-Pentenyl)-3-(1-naphthoyl)indole).
- 424 (IX) JWH-071 (1-Ethyl-3-(1-naphthoyl)indole).
- 425 (X) JWH-072 (1-Propyl-3-(1-naphthoyl)indole).
- 426 (XI) JWH-073 (1-Butyl-3-(1-naphthoyl)indole).
- 427 (XII) JWH-080 (1-Butyl-3-(4-methoxy-1-naphthoyl)indole).
- 428 (XIII) JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl)indole).
- 429 (XIV) JWH-098 (1-Pentyl-2-methyl-3-(4-methoxy-1-
430 naphthoyl)indole).
- 431 (XV) JWH-116 (1-Pentyl-2-ethyl-3-(1-naphthoyl)indole).
- 432 (XVI) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
- 433 (XVII) JWH-149 (1-Pentyl-2-methyl-3-(4-methyl-1-
434 naphthoyl)indole).
- 435 (XVIII) JWH-164 (1-Pentyl-3-(7-methoxy-1-naphthoyl)indole).

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- 436 (XIX) JWH-175 (1-Pentyl-3-(1-naphthylmethyl)indole).
- 437 (XX) JWH-180 (1-Propyl-3-(4-propyl-1-naphthoyl)indole).
- 438 (XXI) JWH-182 (1-Pentyl-3-(4-propyl-1-naphthoyl)indole).
- 439 (XXII) JWH-184 (1-Pentyl-3-[(4-methyl)-1-
- 440 naphthylmethyl]indole).
- 441 (XXIII) JWH-193 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methyl-1-
- 442 naphthoyl)indole).
- 443 (XXIV) JWH-198 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methoxy-1-
- 444 naphthoyl)indole).
- 445 (XXV) JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-
- 446 naphthoyl)indole).
- 447 (XXVI) JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).
- 448 (XXVII) JWH-387 (1-Pentyl-3-(4-bromo-1-naphthoyl)indole).
- 449 (XXVIII) JWH-398 (1-Pentyl-3-(4-chloro-1-naphthoyl)indole).
- 450 (XXIX) JWH-412 (1-Pentyl-3-(4-fluoro-1-naphthoyl)indole).
- 451 (XXX) JWH-424 (1-Pentyl-3-(8-bromo-1-naphthoyl)indole).
- 452 (XXXI) AM-1220 (1-[(1-Methyl-2-piperidinyl)methyl]-3-(1-
- 453 naphthoyl)indole).
- 454 (XXXII) AM-1235 (1-(5-Fluoropentyl)-6-nitro-3-(1-
- 455 naphthoyl)indole).
- 456 (XXXIII) AM-2201 (1-(5-Fluoropentyl)-3-(1-
- 457 naphthoyl)indole).
- 458 (XXXIV) Chloro JWH-018 (1-(Chloropentyl)-3-(1-
- 459 naphthoyl)indole).
- 460 (XXXV) Bromo JWH-018 (1-(Bromopentyl)-3-(1-
- 461 naphthoyl)indole).
- 462 (XXXVI) AM-2232 (1-(4-Cyanobutyl)-3-(1-naphthoyl)indole).
- 463 (XXXVII) THJ-2201 (1-(5-Fluoropentyl)-3-(1-
- 464 naphthoyl)indazole).

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465 (XXXVIII) MAM-2201 (1-(5-Fluoropentyl)-3-(4-methyl-1-
466 naphthoyl)indole).

467 (XXXIX) EAM-2201 (1-(5-Fluoropentyl)-3-(4-ethyl-1-
468 naphthoyl)indole).

469 (XL) EG-018 (9-Pentyl-3-(1-naphthoyl)carbazole).

470 (XLI) EG-2201 (9-(5-Fluoropentyl)-3-(1-
471 naphthoyl)carbazole).

472 c. Naphthoylpyrroles.—Any compound containing a
473 naphthoylpyrrole structure, with or without substitution on the
474 pyrrole ring to any extent, whether or not substituted on the
475 naphthyl ring to any extent, including, but not limited to:

476 (I) JWH-030 (1-Pentyl-3-(1-naphthoyl)pyrrole).

477 (II) JWH-031 (1-Hexyl-3-(1-naphthoyl)pyrrole).

478 (III) JWH-145 (1-Pentyl-5-phenyl-3-(1-naphthoyl)pyrrole).

479 (IV) JWH-146 (1-Heptyl-5-phenyl-3-(1-naphthoyl)pyrrole).

480 (V) JWH-147 (1-Hexyl-5-phenyl-3-(1-naphthoyl)pyrrole).

481 (VI) JWH-307 (1-Pentyl-5-(2-fluorophenyl)-3-(1-
482 naphthoyl)pyrrole).

483 (VII) JWH-309 (1-Pentyl-5-(1-naphthalenyl)-3-(1-
484 naphthoyl)pyrrole).

485 (VIII) JWH-368 (1-Pentyl-5-(3-fluorophenyl)-3-(1-
486 naphthoyl)pyrrole).

487 (IX) JWH-369 (1-Pentyl-5-(2-chlorophenyl)-3-(1-
488 naphthoyl)pyrrole).

489 (X) JWH-370 (1-Pentyl-5-(2-methylphenyl)-3-(1-
490 naphthoyl)pyrrole).

491 d. Naphthylmethylenindenes.—Any compound containing a
492 naphthylmethylenindene structure, with or without substitution
493 at the 3-position of the indene ring to any extent, whether or

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494 not substituted on the naphthyl ring to any extent, including,
495 but not limited to, JWH-176 (3-Pentyl-1-
496 (naphthylmethylene)indene).

497 e. Phenylacetylindoles and Phenylacetylindazoles.—Any
498 compound containing a phenylacetylindole or phenylacetylindazole
499 structure, with or without substitution on the indole or
500 indazole ring to any extent, whether or not substituted on the
501 phenyl ring to any extent, including, but not limited to:

- 502 (I) JWH-167 (1-Pentyl-3-(phenylacetyl)indole).
503 (II) JWH-201 (1-Pentyl-3-(4-methoxyphenylacetyl)indole).
504 (III) JWH-203 (1-Pentyl-3-(2-chlorophenylacetyl)indole).
505 (IV) JWH-250 (1-Pentyl-3-(2-methoxyphenylacetyl)indole).
506 (V) JWH-251 (1-Pentyl-3-(2-methylphenylacetyl)indole).
507 (VI) JWH-302 (1-Pentyl-3-(3-methoxyphenylacetyl)indole).
508 (VII) Cannabipiperidiethanone.
509 (VIII) RCS-8 (1-(2-Cyclohexylethyl)-3-(2-
510 methoxyphenylacetyl)indole).

511 f. Cyclohexylphenols.—Any compound containing a
512 cyclohexylphenol structure, with or without substitution at the
513 5-position of the phenolic ring to any extent, whether or not
514 substituted on the cyclohexyl ring to any extent, including, but
515 not limited to:

- 516 (I) CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-
517 yl)phenol).
518 (II) Cannabicyclohexanol (CP 47,497 dimethyloctyl (C8)
519 homologue).
520 (III) CP-55,940 (2-(3-Hydroxy-6-propanol-cyclohexyl)-5-(2-
521 methyloctan-2-yl)phenol).

522 g. Benzoylindoles and Benzoylindazoles.—Any compound

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523 containing a benzoylindole or benzoylindazole structure, with or
524 without substitution on the indole or indazole ring to any
525 extent, whether or not substituted on the phenyl ring to any
526 extent, including, but not limited to:

527 (I) AM-679 (1-Pentyl-3-(2-iodobenzoyl)indole).

528 (II) AM-694 (1-(5-Fluoropentyl)-3-(2-iodobenzoyl)indole).

529 (III) AM-1241 (1-[(N-Methyl-2-piperidinyl)methyl]-3-(2-
530 iodo-5-nitrobenzoyl)indole).

531 (IV) Pravadoline (1-[2-(4-Morpholinyl)ethyl]-2-methyl-3-(4-
532 methoxybenzoyl)indole).

533 (V) AM-2233 (1-[(N-Methyl-2-piperidinyl)methyl]-3-(2-
534 iodobenzoyl)indole).

535 (VI) RCS-4 (1-Pentyl-3-(4-methoxybenzoyl)indole).

536 (VII) RCS-4 C4 homologue (1-Butyl-3-(4-
537 methoxybenzoyl)indole).

538 (VIII) AM-630 (1-[2-(4-Morpholinyl)ethyl]-2-methyl-6-iodo-
539 3-(4-methoxybenzoyl)indole).

540 h. Tetramethylcyclopropanoylindoles and
541 Tetramethylcyclopropanoylindazoles.—Any compound containing a
542 tetramethylcyclopropanoylindole or
543 tetramethylcyclopropanoylindazole structure, with or without
544 substitution on the indole or indazole ring to any extent,
545 whether or not substituted on the tetramethylcyclopropyl group
546 to any extent, including, but not limited to:

547 (I) UR-144 (1-Pentyl-3-(2,2,3,3-
548 tetramethylcyclopropanoyl)indole).

549 (II) XLR11 (1-(5-Fluoropentyl)-3-(2,2,3,3-
550 tetramethylcyclopropanoyl)indole).

551 (III) Chloro UR-144 (1-(Chloropentyl)-3-(2,2,3,3-

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552 tetramethylcyclopropanoyl) indole).

553 (IV) A-796,260 (1-[2-(4-Morpholinyl)ethyl]-3-(2,2,3,3-
554 tetramethylcyclopropanoyl) indole).

555 (V) A-834,735 (1-[4-(Tetrahydropyranyl)methyl]-3-(2,2,3,3-
556 tetramethylcyclopropanoyl) indole).

557 (VI) M-144 (1-(5-Fluoropentyl)-2-methyl-3-(2,2,3,3-
558 tetramethylcyclopropanoyl) indole).

559 (VII) FUB-144 (1-(4-Fluorobenzyl)-3-(2,2,3,3-
560 tetramethylcyclopropanoyl) indole).

561 (VIII) FAB-144 (1-(5-Fluoropentyl)-3-(2,2,3,3-
562 tetramethylcyclopropanoyl) indazole).

563 (IX) XLR12 (1-(4,4,4-Trifluorobutyl)-3-(2,2,3,3-
564 tetramethylcyclopropanoyl) indole).

565 (X) AB-005 (1-[(1-Methyl-2-piperidinyl)methyl]-3-(2,2,3,3-
566 tetramethylcyclopropanoyl) indole).

567 i. Adamantoylindoles, Adamantoylindazoles, Adamantylindole
568 carboxamides, and Adamantylindazole carboxamides.—Any compound
569 containing an adamantoyl indole, adamantoyl indazole, adamantyl
570 indole carboxamide, or adamantyl indazole carboxamide structure,
571 with or without substitution on the indole or indazole ring to
572 any extent, whether or not substituted on the adamantyl ring to
573 any extent, including, but not limited to:

574 (I) AKB48 (N-Adamant-1-yl 1-pentylindazole-3-carboxamide).

575 (II) Fluoro AKB48 (N-Adamant-1-yl 1-(fluoropentyl)indazole-
576 3-carboxamide).

577 (III) STS-135 (N-Adamant-1-yl 1-(5-fluoropentyl)indole-3-
578 carboxamide).

579 (IV) AM-1248 (1-(1-Methylpiperidine)methyl-3-(1-
580 adamantoyl) indole).

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- 581 (V) AB-001 (1-Pentyl-3-(1-adamantoyl)indole).
- 582 (VI) APICA (N-Adamant-1-yl 1-pentylindole-3-carboxamide).
- 583 (VII) Fluoro AB-001 (1-(Fluoropentyl)-3-(1-
- 584 adamantoyl)indole).
- 585 j. Quinolinyndolecarboxylates,
- 586 Quinolinyndazolecarboxylates, Quinolinyndolecarboxamides,
- 587 and Quinolinyndazolecarboxamides.—Any compound containing a
- 588 quinolinyndole carboxylate, quinolinyndazole carboxylate,
- 589 isoquinolinyndole carboxylate, isoquinolinyndazole
- 590 carboxylate, quinolinyndole carboxamide, quinolinyndazole
- 591 carboxamide, isoquinolinyndole carboxamide, or
- 592 isoquinolinyndazole carboxamide structure, with or without
- 593 substitution on the indole or indazole ring to any extent,
- 594 whether or not substituted on the quinoline or isoquinoline ring
- 595 to any extent, including, but not limited to:
- 596 (I) PB-22 (8-Quinolinyndyl 1-pentylindole-3-carboxylate).
- 597 (II) Fluoro PB-22 (8-Quinolinyndyl 1-(fluoropentyl)indole-3-
- 598 carboxylate).
- 599 (III) BB-22 (8-Quinolinyndyl 1-(cyclohexylmethyl)indole-3-
- 600 carboxylate).
- 601 (IV) FUB-PB-22 (8-Quinolinyndyl 1-(4-fluorobenzyl)indole-3-
- 602 carboxylate).
- 603 (V) NPB-22 (8-Quinolinyndyl 1-pentylindazole-3-carboxylate).
- 604 (VI) Fluoro NPB-22 (8-Quinolinyndyl 1-(fluoropentyl)indazole-
- 605 3-carboxylate).
- 606 (VII) FUB-NPB-22 (8-Quinolinyndyl 1-(4-fluorobenzyl)indazole-
- 607 3-carboxylate).
- 608 (VIII) THJ (8-Quinolinyndyl 1-pentylindazole-3-carboxamide).
- 609 (IX) Fluoro THJ (8-Quinolinyndyl 1-(fluoropentyl)indazole-3-

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610 carboxamide).

611 k. Naphthylindolecarboxylates and
612 Naphthylindazolecarboxylates.—Any compound containing a
613 naphthylindole carboxylate or naphthylindazole carboxylate
614 structure, with or without substitution on the indole or
615 indazole ring to any extent, whether or not substituted on the
616 naphthyl ring to any extent, including, but not limited to:

617 (I) NM-2201 (1-Naphthalenyl 1-(5-fluoropentyl)indole-3-
618 carboxylate).

619 (II) SDB-005 (1-Naphthalenyl 1-pentylindazole-3-
620 carboxylate).

621 (III) Fluoro SDB-005 (1-Naphthalenyl 1-
622 (fluoropentyl)indazole-3-carboxylate).

623 (IV) FDU-PB-22 (1-Naphthalenyl 1-(4-fluorobenzyl)indole-3-
624 carboxylate).

625 (V) 3-CAF (2-Naphthalenyl 1-(2-fluorophenyl)indazole-3-
626 carboxylate).

627 l. Naphthylindole carboxamides and Naphthylindazole
628 carboxamides.—Any compound containing a naphthylindole
629 carboxamide or naphthylindazole carboxamide structure, with or
630 without substitution on the indole or indazole ring to any
631 extent, whether or not substituted on the naphthyl ring to any
632 extent, including, but not limited to:

633 (I) NNEI (N-Naphthalen-1-yl 1-pentylindole-3-carboxamide).

634 (II) Fluoro-NNEI (N-Naphthalen-1-yl 1-(fluoropentyl)indole-
635 3-carboxamide).

636 (III) Chloro-NNEI (N-Naphthalen-1-yl 1-
637 (chloropentyl)indole-3-carboxamide).

638 (IV) MN-18 (N-Naphthalen-1-yl 1-pentylindazole-3-

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639 carboxamide).

640 (V) Fluoro MN-18 (N-Naphthalen-1-yl 1-
641 (fluoropentyl)indazole-3-carboxamide).

642 m. Alkylcarbonyl indole carboxamides, Alkylcarbonyl
643 indazole carboxamides, Alkylcarbonyl indole carboxylates, and
644 Alkylcarbonyl indazole carboxylates.—Any compound containing an
645 alkylcarbonyl group, including 1-amino-3-methyl-1-oxobutan-2-yl,
646 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-amino-1-oxo-3-
647 phenylpropan-2-yl, 1-methoxy-1-oxo-3-phenylpropan-2-yl, with an
648 indole carboxamide, indazole carboxamide, indole carboxylate, or
649 indazole carboxylate, with or without substitution on the indole
650 or indazole ring to any extent, whether or not substituted on
651 the alkylcarbonyl group to any extent, including, but not
652 limited to:

653 (I) ADBICA, (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
654 pentylindole-3-carboxamide).

655 (II) Fluoro ADBICA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-
656 yl)-1-(fluoropentyl)indole-3-carboxamide).

657 (III) Fluoro ABICA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-
658 (fluoropentyl)indole-3-carboxamide).

659 (IV) AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-
660 pentylindazole-3-carboxamide).

661 (V) Fluoro AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-
662 1-(fluoropentyl)indazole-3-carboxamide).

663 (VI) ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-
664 1-pentylindazole-3-carboxamide).

665 (VII) Fluoro ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-
666 oxobutan-2-yl)-1-(fluoropentyl)indazole-3-carboxamide).

667 (VIII) AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-

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668 (4-fluorobenzyl)indazole-3-carboxamide).

669 (IX) ADB-FUBINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-
670 yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).

671 (X) AB-CHMINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-
672 (cyclohexylmethyl)indazole-3-carboxamide).

673 (XI) MA-CHMINACA (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-
674 (cyclohexylmethyl)indazole-3-carboxamide).

675 (XII) MAB-CHMINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-
676 yl)-1-(cyclohexylmethyl)indazole-3-carboxamide).

677 (XIII) AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-
678 pentylindazole-3-carboxamide).

679 (XIV) Fluoro-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-
680 (fluoropentyl)indazole-3-carboxamide).

681 (XV) FUB-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-(4-
682 fluorobenzyl)indazole-3-carboxamide).

683 (XVI) MDMB-CHMINACA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-
684 2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide).

685 (XVII) MDMB-FUBINACA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-
686 2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).

687 (XVIII) MDMB-CHMICA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-
688 2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide).

689 (XIX) PX-1 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(5-
690 fluoropentyl)indole-3-carboxamide).

691 (XX) PX-2 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(5-
692 fluoropentyl)indazole-3-carboxamide).

693 (XXI) PX-3 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-
694 (cyclohexylmethyl)indazole-3-carboxamide).

695 (XXII) PX-4 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(4-
696 fluorobenzyl)indazole-3-carboxamide).

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697 (XXIII) MO-CHMINACA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-
698 2-yl)-1-(cyclohexylmethyl)indazole-3-carboxylate).

699 n. Cumylindolecarboxamides and Cumylindazolecarboxamides.—
700 Any compound containing a N-(2-phenylpropan-2-yl) indole
701 carboxamide or N-(2-phenylpropan-2-yl) indazole carboxamide
702 structure, with or without substitution on the indole or
703 indazole ring to any extent, whether or not substituted on the
704 phenyl ring of the cumyl group to any extent, including, but not
705 limited to:

706 (I) CUMYL-PICA (N-(2-Phenylpropan-2-yl)-1-pentylindole-3-
707 carboxamide).

708 (II) Fluoro CUMYL-PICA (N-(2-Phenylpropan-2-yl)-1-
709 (fluoropentyl)indole-3-carboxamide).

710 o. Other Synthetic Cannabinoids.—Any material, compound,
711 mixture, or preparation that contains any quantity of a
712 Synthetic Cannabinoid, as described in sub-subparagraphs a.-n.:

713 (I) With or without modification or replacement of a
714 carbonyl, carboxamide, alkylene, alkyl, or carboxylate linkage
715 between either two core rings, or linkage between a core ring
716 and group structure, with or without the addition of a carbon or
717 replacement of a carbon;

718 (II) With or without replacement of a core ring or group
719 structure, whether or not substituted on the ring or group
720 structures to any extent; and

721 (III) Is a cannabinoid receptor agonist, unless
722 specifically excepted or unless listed in another schedule or
723 contained within a pharmaceutical product approved by the United
724 States Food and Drug Administration.

725 191. Substituted Cathinones.—Unless specifically excepted,

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726 listed in another schedule, or contained within a pharmaceutical
727 product approved by the United States Food and Drug
728 Administration, any material, compound, mixture, or preparation,
729 including its salts, isomers, esters, or ethers, and salts of
730 isomers, esters, or ethers, whenever the existence of such salts
731 is possible within any of the following specific chemical
732 designations:

733 a. Any compound containing a 2-amino-1-phenyl-1-propanone
734 structure;

735 b. Any compound containing a 2-amino-1-naphthyl-1-propanone
736 structure; or

737 c. Any compound containing a 2-amino-1-thiophenyl-1-
738 propanone structure,

739

740 whether or not the compound is further modified:

741 (I) With or without substitution on the ring system to any
742 extent with alkyl, alkylthio, thio, fused alkylenedioxy, alkoxy,
743 haloalkyl, hydroxyl, nitro, fused furan, fused benzofuran, fused
744 dihydrofuran, fused tetrahydropyran, fused alkyl ring, or halide
745 substituents;

746 (II) With or without substitution at the 3-propanone
747 position with an alkyl substituent or removal of the methyl
748 group at the 3-propanone position;

749 (III) With or without substitution at the 2-amino nitrogen
750 atom with alkyl, dialkyl, acetyl, or benzyl groups, whether or
751 not further substituted in the ring system; or

752 (IV) With or without inclusion of the 2-amino nitrogen atom
753 in a cyclic structure, including, but not limited to:

754 (A) Methcathinone.

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- 755 (B) Ethcathinone.
- 756 (C) Methytlone (3,4-Methylenedioxy~~meth~~cathinone).
- 757 (D) 2,3-Methylenedioxy~~meth~~cathinone.
- 758 (E) MDPV (3,4-Methylenedioxy~~pyro~~valerone).
- 759 (F) Methy~~l~~methcathinone.
- 760 (G) Methoxy~~meth~~cathinone.
- 761 (H) Fluoromethcathinone.
- 762 (I) Methylethcathinone.
- 763 (J) Butylone (3,4-Methylenedioxy-alpha-
- 764 methylaminobutyrophenone).
- 765 (K) Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
- 766 (L) BMDP (3,4-Methylenedioxy-N-benzylcathinone).
- 767 (M) Naphyrone (Naphthylpyrovalerone).
- 768 (N) Bromomethcathinone.
- 769 (O) Buphedrone (alpha-Methylaminobutyrophenone).
- 770 (P) Eutylone (3,4-Methylenedioxy-alpha-
- 771 ethylaminobutyrophenone).
- 772 (Q) Dimethylcathinone.
- 773 (R) Dimethylmethcathinone.
- 774 (S) Pentylone (3,4-Methylenedioxy-alpha-
- 775 methylaminovalerophenone).
- 776 (T) Pentedrone (alpha-Methylaminovalerophenone).
- 777 (U) MDPPP (3,4-Methylenedioxy-alpha-
- 778 pyrrolidinopropiophenone).
- 779 (V) MDPBP (3,4-Methylenedioxy-alpha-
- 780 pyrrolidinobutyrophenone).
- 781 (W) MPPP (Methyl-alpha-pyrrolidinopropiophenone).
- 782 (X) PPP (Pyrrolidinopropiophenone).
- 783 (Y) PVP (Pyrrolidinovalerophenone) or

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784 (Pyrrolidinopentiophenone).

785 (Z) MOPPP (Methoxy-alpha-pyrrolidinopropiophenone).

786 (AA) MPHP (Methyl-alpha-pyrrolidinohexanophenone).

787 (BB) F-MABP (Fluoromethylaminobutyrophenone).

788 (CC) Me-EABP (Methylethylaminobutyrophenone).

789 (DD) PBP (Pyrrolidinobutyrophenone).

790 (EE) MeO-PBP (Methoxypyrrolidinobutyrophenone).

791 (FF) Et-PBP (Ethylpyrrolidinobutyrophenone).

792 (GG) 3-Me-4-MeO-MCAT (3-Methyl-4-Methoxymethcathinone).

793 (HH) Dimethylone (3,4-Methylenedioxy-N,N-

794 dimethylcathinone).

795 (II) 3,4-Methylenedioxy-N,N-diethylcathinone.

796 (JJ) 3,4-Methylenedioxy-N-acetylcathinone.

797 (KK) 3,4-Methylenedioxy-N-acetylmethcathinone.

798 (LL) 3,4-Methylenedioxy-N-acetylethcathinone.

799 (MM) Methylbuphedrone (Methyl-alpha-

800 methylaminobutyrophenone).

801 (NN) Methyl-alpha-methylaminohexanophenone.

802 (OO) N-Ethyl-N-methylcathinone.

803 (PP) PHP (Pyrrolidinohexanophenone).

804 (QQ) PV8 (Pyrrolidinoheptanophenone).

805 (RR) Chloromethcathinone.

806 (SS) 4-Bromo-2,5-dimethoxy-alpha-aminoacetophenone.

807 192. Substituted Phenethylamines.—Unless specifically

808 excepted or unless listed in another schedule, or contained

809 within a pharmaceutical product approved by the United States

810 Food and Drug Administration, any material, compound, mixture,

811 or preparation, including its salts, isomers, esters, or ethers,

812 and salts of isomers, esters, or ethers, whenever the existence

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813 of such salts is possible within any of the following specific
814 chemical designations, any compound containing a phenethylamine
815 structure, without a beta-keto group, and without a benzyl group
816 attached to the amine group, whether or not the compound is
817 further modified with or without substitution on the phenyl ring
818 to any extent with alkyl, alkylthio, nitro, alkoxy, thio,
819 halide, fused alkylenedioxy, fused furan, fused benzofuran,
820 fused dihydrofuran, or fused tetrahydropyran substituents,
821 whether or not further substituted on a ring to any extent, with
822 or without substitution at the alpha or beta position by any
823 alkyl substituent, with or without substitution at the nitrogen
824 atom, and with or without inclusion of the 2-amino nitrogen atom
825 in a cyclic structure, including, but not limited to:

- 826 a. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
827 b. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
828 c. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
829 d. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
830 e. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
831 f. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
832 g. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
833 h. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
834 i. 2C-D (4-Methyl-2,5-dimethoxyphenethylamine).
835 j. 2C-H (2,5-Dimethoxyphenethylamine).
836 k. 2C-N (4-Nitro-2,5-dimethoxyphenethylamine).
837 l. 2C-P (4-(n)-Propyl-2,5-dimethoxyphenethylamine).
838 m. MDMA (3,4-Methylenedioxyamphetamin).
839 n. MBDB (Methylbenzodioxolylbutanamine) or (3,4-
840 Methylenedioxy-N-methylbutanamine).
841 o. MDA (3,4-Methylenedioxyamphetamin).

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- 842 p. 2,5-Dimethoxyamphetamine.
- 843 q. Fluoroamphetamine.
- 844 r. Fluoromethamphetamine.
- 845 s. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
- 846 t. DOB (4-Bromo-2,5-dimethoxyamphetamine).
- 847 u. DOC (4-Chloro-2,5-dimethoxyamphetamine).
- 848 v. DOET (4-Ethyl-2,5-dimethoxyamphetamine).
- 849 w. DOI (4-Iodo-2,5-dimethoxyamphetamine).
- 850 x. DOM (4-Methyl-2,5-dimethoxyamphetamine).
- 851 y. PMA (4-Methoxyamphetamine).
- 852 z. N-Ethylamphetamine.
- 853 aa. 3,4-Methylenedioxy-N-hydroxyamphetamine.
- 854 bb. 5-Methoxy-3,4-methylenedioxyamphetamine.
- 855 cc. PMMA (4-Methoxymethamphetamine).
- 856 dd. N,N-Dimethylamphetamine.
- 857 ee. 3,4,5-Trimethoxyamphetamine.
- 858 ff. 4-APB (4-(2-Aminopropyl)benzofuran).
- 859 gg. 5-APB (5-(2-Aminopropyl)benzofuran).
- 860 hh. 6-APB (6-(2-Aminopropyl)benzofuran).
- 861 ii. 7-APB (7-(2-Aminopropyl)benzofuran).
- 862 jj. 4-APDB (4-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 863 kk. 5-APDB (5-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 864 ll. 6-APDB (6-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 865 mm. 7-APDB (7-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 866 nn. 4-MAPB (4-(2-Methylaminopropyl)benzofuran).
- 867 oo. 5-MAPB (5-(2-Methylaminopropyl)benzofuran).
- 868 pp. 6-MAPB (6-(2-Methylaminopropyl)benzofuran).
- 869 qq. 7-MAPB (7-(2-Methylaminopropyl)benzofuran).
- 870 rr. 5-EAPB (5-(2-Ethylaminopropyl)benzofuran).

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871 ss. 5-MAPDB (5-(2-Methylaminopropyl)-2,3-
872 dihydrobenzofuran),

873

874 which does not include phenethylamine, mescaline as described in
875 subparagraph 20., substituted cathinones as described in
876 subparagraph 191., N-Benzyl phenethylamine compounds as
877 described in subparagraph 193., or methamphetamine as described
878 in subparagraph (2)(c)5.

879 193. N-Benzyl Phenethylamine Compounds.—Unless specifically
880 excepted or unless listed in another schedule, or contained
881 within a pharmaceutical product approved by the United States
882 Food and Drug Administration, any material, compound, mixture,
883 or preparation, including its salts, isomers, esters, or ethers,
884 and salts of isomers, esters, or ethers, whenever the existence
885 of such salts is possible within any of the following specific
886 chemical designations, any compound containing a phenethylamine
887 structure without a beta-keto group, with substitution on the
888 nitrogen atom of the amino group with a benzyl substituent, with
889 or without substitution on the phenyl or benzyl ring to any
890 extent with alkyl, alkoxy, thio, alkylthio, halide, fused
891 alkylendioxy, fused furan, fused benzofuran, or fused
892 tetrahydropyran substituents, whether or not further substituted
893 on a ring to any extent, with or without substitution at the
894 alpha position by any alkyl substituent, including, but not
895 limited to:

896 a. 25B-NBOMe (4-Bromo-2,5-dimethoxy-[N-(2-
897 methoxybenzyl)]phenethylamine).

898 b. 25B-NBOH (4-Bromo-2,5-dimethoxy-[N-(2-
899 hydroxybenzyl)]phenethylamine).

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- 900 c. 25B-NBF (4-Bromo-2,5-dimethoxy-[N-(2-
901 fluorobenzyl)]phenethylamine).
- 902 d. 25B-NBMD (4-Bromo-2,5-dimethoxy-[N-(2,3-
903 methylenedioxybenzyl)]phenethylamine).
- 904 e. 25I-NBOMe (4-Iodo-2,5-dimethoxy-[N-(2-
905 methoxybenzyl)]phenethylamine).
- 906 f. 25I-NBOH (4-Iodo-2,5-dimethoxy-[N-(2-
907 hydroxybenzyl)]phenethylamine).
- 908 g. 25I-NBF (4-Iodo-2,5-dimethoxy-[N-(2-
909 fluorobenzyl)]phenethylamine).
- 910 h. 25I-NBMD (4-Iodo-2,5-dimethoxy-[N-(2,3-
911 methylenedioxybenzyl)]phenethylamine).
- 912 i. 25T2-NBOMe (4-Methylthio-2,5-dimethoxy-[N-(2-
913 methoxybenzyl)]phenethylamine).
- 914 j. 25T4-NBOMe (4-Isopropylthio-2,5-dimethoxy-[N-(2-
915 methoxybenzyl)]phenethylamine).
- 916 k. 25T7-NBOMe (4-(n)-Propylthio-2,5-dimethoxy-[N-(2-
917 methoxybenzyl)]phenethylamine).
- 918 l. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-
919 methoxybenzyl)]phenethylamine).
- 920 m. 25C-NBOH (4-Chloro-2,5-dimethoxy-[N-(2-
921 hydroxybenzyl)]phenethylamine).
- 922 n. 25C-NBF (4-Chloro-2,5-dimethoxy-[N-(2-
923 fluorobenzyl)]phenethylamine).
- 924 o. 25C-NBMD (4-Chloro-2,5-dimethoxy-[N-(2,3-
925 methylenedioxybenzyl)]phenethylamine).
- 926 p. 25H-NBOMe (2,5-Dimethoxy-[N-(2-
927 methoxybenzyl)]phenethylamine).
- 928 q. 25H-NBOH (2,5-Dimethoxy-[N-(2-

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929 hydroxybenzyl)]phenethylamine).

930 r. 25H-NBF (2,5-Dimethoxy-[N-(2-
931 fluorobenzyl)]phenethylamine).

932 s. 25D-NBOMe (4-Methyl-2,5-dimethoxy-[N-(2-
933 methoxybenzyl)]phenethylamine),

934

935 which does not include substituted cathinones as described in
936 subparagraph 191.

937 194. Substituted Tryptamines.—Unless specifically excepted
938 or unless listed in another schedule, or contained within a
939 pharmaceutical product approved by the United States Food and
940 Drug Administration, any material, compound, mixture, or
941 preparation containing a 2-(1H-indol-3-yl)ethanamine, for
942 example tryptamine, structure with or without mono- or di-
943 substitution of the amine nitrogen with alkyl or alkenyl groups,
944 or by inclusion of the amino nitrogen atom in a cyclic
945 structure, whether or not substituted at the alpha position with
946 an alkyl group, whether or not substituted on the indole ring to
947 any extent with any alkyl, alkoxy, halo, hydroxyl, or acetoxy
948 groups, including, but not limited to:

949 a. Alpha-Ethyltryptamine.

950 b. Bufotenine.

951 c. DET (Diethyltryptamine).

952 d. DMT (Dimethyltryptamine).

953 e. MET (N-Methyl-N-ethyltryptamine).

954 f. DALT (N,N-Diallyltryptamine).

955 g. EiPT (N-Ethyl-N-isopropyltryptamine).

956 h. MiPT (N-Methyl-N-isopropyltryptamine).

957 i. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).

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- 958 j. 5-Hydroxy-N-methyltryptamine.
- 959 k. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).
- 960 l. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
- 961 m. Methyltryptamine.
- 962 n. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
- 963 o. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
- 964 p. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
- 965 q. DiPT (N,N-Diisopropyltryptamine).
- 966 r. DPT (N,N-Dipropyltryptamine).
- 967 s. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
- 968 t. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
- 969 u. 4-AcO-DMT (4-Acetoxy-N,N-dimethyltryptamine).
- 970 v. 4-AcO-DiPT (4-Acetoxy-N,N-diisopropyltryptamine).
- 971 w. 4-Hydroxy-DET (4-Hydroxy-N,N-diethyltryptamine).
- 972 x. 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-ethyltryptamine).
- 973 y. 4-Hydroxy-MiPT (4-Hydroxy-N-methyl-N-
- 974 isopropyltryptamine).
- 975 z. Methyl-alpha-ethyltryptamine.
- 976 aa. Bromo-DALT (Bromo-N,N-diallyltryptamine),
- 977

978 which does not include tryptamine, psilocyn as described in
 979 subparagraph 34., or psilocybin as described in subparagraph 33.

980 195. Substituted Phenylcyclohexylamines.—Unless
 981 specifically excepted or unless listed in another schedule, or
 982 contained within a pharmaceutical product approved by the United
 983 States Food and Drug Administration, any material, compound,
 984 mixture, or preparation containing a phenylcyclohexylamine
 985 structure, with or without any substitution on the phenyl ring,
 986 any substitution on the cyclohexyl ring, any replacement of the

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- 987 phenyl ring with a thiophenyl or benzothiophenyl ring, with or
988 without substitution on the amine with alkyl, dialkyl, or alkoxy
989 substituents, inclusion of the nitrogen in a cyclic structure,
990 or any combination of the above, including, but not limited to:
- 991 a. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP
992 (Benocyclidine).
 - 993 b. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine analog
994 of phencyclidine).
 - 995 c. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine
996 analog of phencyclidine).
 - 997 d. PCPr (Phenylcyclohexylpropylamine).
 - 998 e. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine) (Thiophene
999 analog of phencyclidine).
 - 1000 f. PCEEA (Phenylcyclohexyl(ethoxyethylamine)).
 - 1001 g. PCMPA (Phenylcyclohexyl(methoxypropylamine)).
 - 1002 h. Methoxetamine.
 - 1003 i. 3-Methoxy-PCE ((3-Methoxyphenyl)cyclohexylethylamine).
 - 1004 j. Bromo-PCP ((Bromophenyl)cyclohexylpiperidine).
 - 1005 k. Chloro-PCP ((Chlorophenyl)cyclohexylpiperidine).
 - 1006 l. Fluoro-PCP ((Fluorophenyl)cyclohexylpiperidine).
 - 1007 m. Hydroxy-PCP ((Hydroxyphenyl)cyclohexylpiperidine).
 - 1008 n. Methoxy-PCP ((Methoxyphenyl)cyclohexylpiperidine).
 - 1009 o. Methyl-PCP ((Methylphenyl)cyclohexylpiperidine).
 - 1010 p. Nitro-PCP ((Nitrophenyl)cyclohexylpiperidine).
 - 1011 q. Oxo-PCP ((Oxophenyl)cyclohexylpiperidine).
 - 1012 r. Amino-PCP ((Aminophenyl)cyclohexylpiperidine).
 - 1013 196. W-15, 4-chloro-N-[1-(2-phenylethyl)-2-
1014 piperidinylidene]-benzenesulfonamide.
 - 1015 197. W-18, 4-chloro-N-[1-[2-(4-nitrophenyl)ethyl]-2-

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1016 piperidinylidene]-benzenesulfonamide.

1017 198. AH-7921, 3,4-dichloro-N-[[1-
1018 (dimethylamino)cyclohexyl]methyl]-benzamide.

1019 199. U47700, trans-3,4-dichloro-N-[2-
1020 (dimethylamino)cyclohexyl]-N-methyl-benzamide.

1021 200. MT-45, 1-cyclohexyl-4-(1,2-diphenylethyl)-piperazine,
1022 dihydrochloride.

1023 Section 4. Paragraph (i) of subsection (1) of section
1024 893.13, Florida Statutes, is amended to read:

1025 893.13 Prohibited acts; penalties.—

1026 (1)

1027 (i) Except as authorized by this chapter, a person commits
1028 a felony of the first degree, punishable as provided in s.
1029 775.082, s. 775.083, or s. 775.084, and must be sentenced to a
1030 mandatory minimum term of imprisonment of 3 years, if:

1031 1. The person sells, manufactures, or delivers, or
1032 possesses with intent to sell, manufacture, or deliver, any of
1033 the following:

1034 a. Alfentanil, as described in s. 893.03(2)(b)1.;

1035 b. Carfentanil, as described in s. 893.03(2)(b)6.;

1036 c. Fentanyl, as described in s. 893.03(2)(b)9.;

1037 d. Sufentanil, as described in s. 893.03(2)(b)30.;

1038 e. A fentanyl derivative, as described in s.
1039 893.03(1)(a)63.;

1040 f. Xylazine, as described in s. 893.03(1)(c)37.;

1041 g.f. A controlled substance analog, as described in s.
1042 893.0356, of any substance described in sub-subparagraphs a.-f.
1043 ~~sub-subparagraphs a.-e.~~; or

1044 h.g. A mixture containing any substance described in sub-

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1045 subparagraphs a.-g. ~~sub-subparagraphs a.-f.~~; and

1046 2. The substance or mixture listed in subparagraph 1. is in
1047 a form that resembles, or is mixed, granulated, absorbed, spray-
1048 dried, or aerosolized as or onto, coated on, in whole or in
1049 part, or solubilized with or into, a product, when such product
1050 or its packaging further has at least one of the following
1051 attributes:

1052 a. Resembles the trade dress of a branded food product,
1053 consumer food product, or logo food product;

1054 b. Incorporates an actual or fake registered copyright,
1055 service mark, or trademark;

1056 c. Resembles candy, cereal, a gummy, a vitamin, or a
1057 chewable product, such as a gum or gelatin-based product; or

1058 d. Contains a cartoon character imprint.

1059 Section 5. Paragraph (c) of subsection (1) of section
1060 893.135, Florida Statutes, is amended to read:

1061 893.135 Trafficking; mandatory sentences; suspension or
1062 reduction of sentences; conspiracy to engage in trafficking.-

1063 (1) Except as authorized in this chapter or in chapter 499
1064 and notwithstanding the provisions of s. 893.13:

1065 (c)1. A person who knowingly sells, purchases,
1066 manufactures, delivers, or brings into this state, or who is
1067 knowingly in actual or constructive possession of, 4 grams or
1068 more of any morphine, opium, hydromorphone, or any salt,
1069 derivative, isomer, or salt of an isomer thereof, including
1070 heroin, as described in s. 893.03(1)(b), (2)(a), (3)(c)3., or
1071 (3)(c)4., or 4 grams or more of any mixture containing any such
1072 substance, but less than 30 kilograms of such substance or
1073 mixture, commits a felony of the first degree, which felony

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1074 shall be known as "trafficking in illegal drugs," punishable as
1075 provided in s. 775.082, s. 775.083, or s. 775.084. If the
1076 quantity involved:

1077 a. Is 4 grams or more, but less than 14 grams, such person
1078 shall be sentenced to a mandatory minimum term of imprisonment
1079 of 3 years and shall be ordered to pay a fine of \$50,000.

1080 b. Is 14 grams or more, but less than 28 grams, such person
1081 shall be sentenced to a mandatory minimum term of imprisonment
1082 of 15 years and shall be ordered to pay a fine of \$100,000.

1083 c. Is 28 grams or more, but less than 30 kilograms, such
1084 person shall be sentenced to a mandatory minimum term of
1085 imprisonment of 25 years and shall be ordered to pay a fine of
1086 \$500,000.

1087 2. A person who knowingly sells, purchases, manufactures,
1088 delivers, or brings into this state, or who is knowingly in
1089 actual or constructive possession of, 28 grams or more of
1090 hydrocodone, as described in s. 893.03(2)(a)1.k., codeine, as
1091 described in s. 893.03(2)(a)1.g., or any salt thereof, or 28
1092 grams or more of any mixture containing any such substance,
1093 commits a felony of the first degree, which felony shall be
1094 known as "trafficking in hydrocodone," punishable as provided in
1095 s. 775.082, s. 775.083, or s. 775.084. If the quantity involved:

1096 a. Is 28 grams or more, but less than 50 grams, such person
1097 shall be sentenced to a mandatory minimum term of imprisonment
1098 of 3 years and shall be ordered to pay a fine of \$50,000.

1099 b. Is 50 grams or more, but less than 100 grams, such
1100 person shall be sentenced to a mandatory minimum term of
1101 imprisonment of 7 years and shall be ordered to pay a fine of
1102 \$100,000.

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1103 c. Is 100 grams or more, but less than 300 grams, such
1104 person shall be sentenced to a mandatory minimum term of
1105 imprisonment of 15 years and shall be ordered to pay a fine of
1106 \$500,000.

1107 d. Is 300 grams or more, but less than 30 kilograms, such
1108 person shall be sentenced to a mandatory minimum term of
1109 imprisonment of 25 years and shall be ordered to pay a fine of
1110 \$750,000.

1111 3. A person who knowingly sells, purchases, manufactures,
1112 delivers, or brings into this state, or who is knowingly in
1113 actual or constructive possession of, 7 grams or more of
1114 oxycodone, as described in s. 893.03(2)(a)1.q., or any salt
1115 thereof, or 7 grams or more of any mixture containing any such
1116 substance, commits a felony of the first degree, which felony
1117 shall be known as "trafficking in oxycodone," punishable as
1118 provided in s. 775.082, s. 775.083, or s. 775.084. If the
1119 quantity involved:

1120 a. Is 7 grams or more, but less than 14 grams, such person
1121 shall be sentenced to a mandatory minimum term of imprisonment
1122 of 3 years and shall be ordered to pay a fine of \$50,000.

1123 b. Is 14 grams or more, but less than 25 grams, such person
1124 shall be sentenced to a mandatory minimum term of imprisonment
1125 of 7 years and shall be ordered to pay a fine of \$100,000.

1126 c. Is 25 grams or more, but less than 100 grams, such
1127 person shall be sentenced to a mandatory minimum term of
1128 imprisonment of 15 years and shall be ordered to pay a fine of
1129 \$500,000.

1130 d. Is 100 grams or more, but less than 30 kilograms, such
1131 person shall be sentenced to a mandatory minimum term of

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1132 imprisonment of 25 years and shall be ordered to pay a fine of
1133 \$750,000.

1134 4.a. A person who knowingly sells, purchases, manufactures,
1135 delivers, or brings into this state, or who is knowingly in
1136 actual or constructive possession of, 4 grams or more of:

1137 (I) Alfentanil, as described in s. 893.03(2)(b)1.;

1138 (II) Carfentanil, as described in s. 893.03(2)(b)6.;

1139 (III) Fentanyl, as described in s. 893.03(2)(b)9.;

1140 (IV) Sufentanil, as described in s. 893.03(2)(b)30.;

1141 (V) A fentanyl derivative, as described in s.

1142 893.03(1)(a)63.;

1143 (VI) A controlled substance analog, as described in s.
1144 893.0356, of any substance described in sub-sub-subparagraphs
1145 (I)-(V); or

1146 (VII) A mixture containing any substance described in sub-
1147 sub-subparagraphs (I)-(VI),

1148

1149 commits a felony of the first degree, which felony shall be
1150 known as "trafficking in dangerous fentanyl or fentanyl
1151 analogues," punishable as provided in s. 775.082, s. 775.083, or
1152 s. 775.084.

1153 b. If the quantity involved under sub-subparagraph a.:

1154 (I) Is 4 grams or more, but less than 14 grams, such person
1155 shall be sentenced to a mandatory minimum term of imprisonment
1156 of 7 years~~7~~ and shall be ordered to pay a fine of \$50,000.

1157 (II) Is 14 grams or more, but less than 28 grams, such
1158 person shall be sentenced to a mandatory minimum term of
1159 imprisonment of 20 years~~7~~ and shall be ordered to pay a fine of
1160 \$100,000.

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1161 (III) Is 28 grams or more, such person shall be sentenced
1162 to a mandatory minimum term of imprisonment of 25 years~~,~~ and
1163 shall be ordered to pay a fine of \$500,000.

1164 c. A person 18 years of age or older who violates sub-
1165 subparagraph a. by knowingly selling or delivering to a minor at
1166 least 4 grams of a substance or mixture listed in sub-
1167 subparagraph a. shall be sentenced to a mandatory minimum term
1168 of not less than 25 years and not exceeding life imprisonment,
1169 and shall be ordered to pay a fine of \$1 million if the
1170 substance or mixture listed in sub-subparagraph a. is in a form
1171 that resembles, or is mixed, granulated, absorbed, spray-dried,
1172 or aerosolized as or onto, coated on, in whole or in part, or
1173 solubilized with or into, a product, when such product or its
1174 packaging further has at least one of the following attributes:

1175 (I) Resembles the trade dress of a branded food product,
1176 consumer food product, or logo food product;

1177 (II) Incorporates an actual or fake registered copyright,
1178 service mark, or trademark;

1179 (III) Resembles candy, cereal, a gummy, a vitamin, or a
1180 chewable product, such as a gum or gelatin-based product; or

1181 (IV) Contains a cartoon character imprint.

1182 5. A person who knowingly sells, purchases, manufactures,
1183 delivers, or brings into this state, or who is knowingly in
1184 actual or constructive possession of, 30 kilograms or more of
1185 any morphine, opium, oxycodone, hydrocodone, codeine,
1186 hydromorphone, or any salt, derivative, isomer, or salt of an
1187 isomer thereof, including heroin, as described in s.
1188 893.03(1)(b), (2)(a), (3)(c)3., or (3)(c)4., or 30 kilograms or
1189 more of any mixture containing any such substance, commits the

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1190 first degree felony of trafficking in illegal drugs. A person
1191 who has been convicted of the first degree felony of trafficking
1192 in illegal drugs under this subparagraph shall be punished by
1193 life imprisonment and is ineligible for any form of
1194 discretionary early release except pardon or executive clemency
1195 or conditional medical release under s. 947.149. However, if the
1196 court determines that, in addition to committing any act
1197 specified in this paragraph:

1198 a. The person intentionally killed an individual or
1199 counseled, commanded, induced, procured, or caused the
1200 intentional killing of an individual and such killing was the
1201 result; or

1202 b. The person's conduct in committing that act led to a
1203 natural, though not inevitable, lethal result,

1204

1205 such person commits the capital felony of trafficking in illegal
1206 drugs, punishable as provided in ss. 775.082 and 921.142. A
1207 person sentenced for a capital felony under this paragraph shall
1208 also be sentenced to pay the maximum fine provided under
1209 subparagraph 1.

1210 6. A person who knowingly brings into this state 60
1211 kilograms or more of any morphine, opium, oxycodone,
1212 hydrocodone, codeine, hydromorphone, or any salt, derivative,
1213 isomer, or salt of an isomer thereof, including heroin, as
1214 described in s. 893.03(1)(b), (2)(a), (3)(c)3., or (3)(c)4., or
1215 60 kilograms or more of any mixture containing any such
1216 substance, and who knows that the probable result of such
1217 importation would be the death of a person, commits capital
1218 importation of illegal drugs, a capital felony punishable as

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1219 provided in ss. 775.082 and 921.142. A person sentenced for a
1220 capital felony under this paragraph shall also be sentenced to
1221 pay the maximum fine provided under subparagraph 1.

1222 7. A person who knowingly sells, purchases, manufactures,
1223 delivers, or brings into this state, or who is knowingly in
1224 actual or constructive possession of, 28 grams or more of
1225 xylazine, as described in s. 893.03(1)(c)37., or any salt
1226 thereof, or 28 grams or more of any mixture containing any such
1227 substance, commits a felony of the first degree, which felony
1228 shall be known as "trafficking in xylazine," punishable as
1229 provided in s. 775.082, s. 775.083, or s. 775.084. If the
1230 quantity involved:

1231 a. Is 28 grams or more, but less than 100 grams, such
1232 person shall be sentenced to a mandatory minimum term of
1233 imprisonment of 3 years and shall be ordered to pay a fine of
1234 \$50,000.

1235 b. Is 100 grams or more, but less than 200 grams, such
1236 person shall be sentenced to a mandatory minimum term of
1237 imprisonment of 7 years and shall be ordered to pay a fine of
1238 \$100,000.

1239 c. Is 200 grams or more, such person shall be sentenced to
1240 a mandatory minimum term of imprisonment of 25 years and shall
1241 be ordered to pay a fine of \$500,000.

1242 Section 6. Except as otherwise expressly provided in this
1243 act and except for this section, which shall take effect upon
1244 this act becoming a law, this act shall take effect October 1,
1245 2026.