

By Senator Yarborough

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A bill to be entitled
An act relating to controlled substances; amending s.
893.03, F.S.; adding 7-Hydroxymitragynine as a
Schedule I controlled substance; excepting from the
list of Schedule I controlled substances certain
xylazine animal drug products approved by the United
States Food and Drug Administration and used for
certain purposes; amending s. 893.13, F.S.; providing
criminal penalties and requiring a mandatory minimum
term of imprisonment if a person sells, manufactures,
or delivers or possesses with intent to sell,
manufacture, or deliver xylazine; amending s. 893.135,
F.S.; creating the offense of trafficking in xylazine;
providing criminal penalties and requiring a mandatory
minimum term of imprisonment and fines based on the
quantity of the controlled substance involved in the
offense; providing effective dates.

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

Section 1. Effective July 1, 2026, paragraphs (a) and (c)
of subsection (1) of section 893.03, Florida Statutes, are
amended to read:

893.03 Standards and schedules.—The substances enumerated
in this section are controlled by this chapter. The controlled
substances listed or to be listed in Schedules I, II, III, IV,
and V are included by whatever official, common, usual,
chemical, trade name, or class designated. The provisions of
this section shall not be construed to include within any of the

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schedules contained in this section any excluded drugs listed within the purview of 21 C.F.R. s. 1308.22, styled "Excluded Substances"; 21 C.F.R. s. 1308.24, styled "Exempt Chemical Preparations"; 21 C.F.R. s. 1308.32, styled "Exempted Prescription Products"; or 21 C.F.R. s. 1308.34, styled "Exempt Anabolic Steroid Products."

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

(a) Unless specifically excepted or unless listed in another schedule, any of the following substances, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation:

1. Acetyl-alpha-methylfentanyl.
2. Acetylmethadol.
3. Allylprodine.
4. Alphacetylmethadol (except levo-alphacetylmethadol, also known as levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM).
5. Alphamethadol.
6. Alpha-methylfentanyl (N-[1-(alpha-methyl-betaphenyl) ethyl-4-piperidyl] propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine).
7. Alpha-methylthiofentanyl.
8. Alphameprodine.

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59 9. Benzethidine.
60 10. Benzylfentanyl.
61 11. Betacetylmethadol.
62 12. Beta-hydroxyfentanyl.
63 13. Beta-hydroxy-3-methylfentanyl.
64 14. Betameprodine.
65 15. Betamethadol.
66 16. Betaprodine.
67 17. Clonitazene.
68 18. Dextromoramide.
69 19. Diampromide.
70 20. Diethylthiambutene.
71 21. Difenoxin.
72 22. Dimenoxadol.
73 23. Dimepheptanol.
74 24. Dimethylthiambutene.
75 25. Dioxaphetyl butyrate.
76 26. Dipipanone.
77 27. Ethylmethylthiambutene.
78 28. Etonitazene.
79 29. Etoxeridine.
80 30. Flunitrazepam.
81 31. Furethidine.
82 32. Hydroxypethidine.
83 33. Ketobemidone.
84 34. Levomoramide.
85 35. Levophenacylmorphane.
86 36. Desmethyprodine (1-Methyl-4-Phenyl-4-
87 Propionoxypiperidine).

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37. 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide).

38. 3-Methylthiofentanyl.

39. Morpheridine.

40. Noracymethadol.

41. Norlevorphanol.

42. Normethadone.

43. Norpipanone.

44. Para-Fluorofentanyl.

45. Phenadoxone.

46. Phenampromide.

47. Phenomorphan.

48. Phenoperidine.

49. PEPAP (1-(2-Phenylethyl)-4-Phenyl-4-Acetyloxypiperidine).

50. Piritramide.

51. Proheptazine.

52. Properidine.

53. Propiram.

54. Racemoramide.

55. Thenylfentanyl.

56. Thiofentanyl.

57. Tianeptine.

58. Tilidine.

59. Trimeperidine.

60. Acetylfentanyl.

61. Butyrylfentanyl.

62. Beta-Hydroxythiofentanyl.

63. Fentanyl derivatives. Unless specifically excepted,

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117 listed in another schedule, or contained within a pharmaceutical
118 product approved by the United States Food and Drug
119 Administration, any material, compound, mixture, or preparation,
120 including its salts, isomers, esters, or ethers, and salts of
121 isomers, esters, or ethers, whenever the existence of such salts
122 is possible within any of the following specific chemical
123 designations containing a 4-anilidopiperidine structure:

124 a. With or without substitution at the carbonyl of the
125 aniline moiety with alkyl, alkenyl, carboalkoxy, cycloalkyl,
126 methoxyalkyl, cyanoalkyl, or aryl groups, or furanyl,
127 dihydrofuranyl, benzyl moiety, or rings containing heteroatoms
128 sulfur, oxygen, or nitrogen;

129 b. With or without substitution at the piperidine amino
130 moiety with a phenethyl, benzyl, alkylaryl (including
131 heteroaromatics), alkyltetrazolyl ring, or an alkyl or
132 carbomethoxy group, whether or not further substituted in the
133 ring or group;

134 c. With or without substitution or addition to the
135 piperidine ring to any extent with one or more methyl,
136 carbomethoxy, methoxy, methoxymethyl, aryl, allyl, or ester
137 groups;

138 d. With or without substitution of one or more hydrogen
139 atoms for halogens, or methyl, alkyl, or methoxy groups, in the
140 aromatic ring of the anilide moiety;

141 e. With or without substitution at the alpha or beta
142 position of the piperidine ring with alkyl, hydroxyl, or methoxy
143 groups;

144 f. With or without substitution of the benzene ring of the
145 anilide moiety for an aromatic heterocycle; and

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g. With or without substitution of the piperidine ring for a pyrrolidine ring, perhydroazepine ring, or azepine ring;

excluding, Alfentanil, Carfentanil, Fentanyl, and Sufentanil; including, but not limited to:

(I) Acetyl-alpha-methylfentanyl.

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

(III) Alpha-methylthiofentanyl.

(IV) Benzylfentanyl.

(V) Beta-hydroxyfentanyl.

(VI) Beta-hydroxy-3-methylfentanyl.

(VII) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide).

(VIII) 3-Methylthiofentanyl.

(IX) Para-Fluorofentanyl.

(X) Thenylfentanyl or Thienyl fentanyl.

(XI) Thiofentanyl.

(XII) Acetylfentanyl.

(XIII) Butyrylfentanyl.

(XIV) Beta-Hydroxythiofentanyl.

(XV) Lofentanil.

(XVI) Ocfentanil.

(XVII) Ohmfentanyl.

(XVIII) Benzodioxolefentanyl.

(XIX) Furanyl fentanyl.

(XX) Pentanoyl fentanyl.

(XXI) Cyclopentyl fentanyl.

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(XXII) Isobutyryl fentanyl.

(XXIII) Remifentanil.

64. Nitazene derivatives. Unless specifically excepted, listed in another schedule, or contained within a pharmaceutical product approved by the United States Food and Drug Administration, any material, compound, mixture, or preparation, including its salts, isomers, esters, or ethers, and salts of isomers, esters, or ethers, whenever the existence of such salts is possible within any of the following specific chemical designations containing a benzimidazole ring with an ethylamine substitution at the 1-position and a benzyl ring substitution at the 2-position structure:

a. With or without substitution on the benzimidazole ring with alkyl, alkoxy, carboalkoxy, amino, nitro, or aryl groups, or halogens;

b. With or without substitution at the ethylamine amino moiety with alkyl, dialkyl, acetyl, or benzyl groups, whether or not further substituted in the ring system;

c. With or without inclusion of the ethylamine amino moiety in a cyclic structure;

d. With or without substitution of the benzyl ring; or

e. With or without replacement of the benzyl ring with an aromatic ring, including, but not limited to:

(I) Butonitazene.

(II) Clonitazene.

(III) Etodesnitazene.

(IV) Etonitazene.

(V) Flunitazene.

(VI) Isotodesnitazene.

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(VII) Isotonitazene.

(VIII) Metodesnitazene.

(IX) Metonitazene.

(X) Nitazene.

(XI) N-Desethyl Etonitazene.

(XII) N-Desethyl Isotonitazene.

(XIII) N-Piperidino Etonitazene.

(XIV) N-Pyrrolidino Etonitazene.

(XV) Protonitazene.

65. 7-Hydroxymitragynine (methyl (E)-2-[(2S,3S,7aS,12bS)-3-ethyl-7a-hydroxy-8-methoxy-2,3,4,6,7,12b-hexahydro-1H-indolo[2,3-a]quinolizin-2-yl]-3-methoxyprop-2-enoate) concentrated at a level above 400 parts per million on a dry-weight basis.

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

1. Alpha-Ethyltryptamine.

2. 4-Methylaminorex (2-Amino-4-methyl-5-phenyl-2-oxazoline).

3. Aminorex (2-Amino-5-phenyl-2-oxazoline).

4. DOB (4-Bromo-2,5-dimethoxyamphetamine).

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- 233 5. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
- 234 6. Bufotenine.
- 235 7. Cannabis.
- 236 8. Cathinone.
- 237 9. DET (Diethyltryptamine).
- 238 10. 2,5-Dimethoxyamphetamine.
- 239 11. DOET (4-Ethyl-2,5-Dimethoxyamphetamine).
- 240 12. DMT (Dimethyltryptamine).
- 241 13. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine
- 242 analog of phencyclidine).
- 243 14. JB-318 (N-Ethyl-3-piperidyl benzilate).
- 244 15. N-Ethylamphetamine.
- 245 16. Fenethylamine.
- 246 17. 3,4-Methylenedioxy-N-hydroxyamphetamine.
- 247 18. Ibogaine.
- 248 19. LSD (Lysergic acid diethylamide).
- 249 20. Mescaline.
- 250 21. Methcathinone.
- 251 22. 5-Methoxy-3,4-methylenedioxyamphetamine.
- 252 23. PMA (4-Methoxyamphetamine).
- 253 24. PMMA (4-Methoxymethamphetamine).
- 254 25. DOM (4-Methyl-2,5-dimethoxyamphetamine).
- 255 26. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
- 256 27. MDA (3,4-Methylenedioxyamphetamine).
- 257 28. JB-336 (N-Methyl-3-piperidyl benzilate).
- 258 29. N,N-Dimethylamphetamine.
- 259 30. Parahexyl.
- 260 31. Peyote.
- 261 32. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine

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262 analog of phencyclidine).

263 33. Psilocybin.

264 34. Psilocyn.

265 35. *Salvia divinorum*, except for any drug product approved
266 by the United States Food and Drug Administration which contains
267 *Salvia divinorum* or its isomers, esters, ethers, salts, and
268 salts of isomers, esters, and ethers, if the existence of such
269 isomers, esters, ethers, and salts is possible within the
270 specific chemical designation.

271 36. Salvinorin A, except for any drug product approved by
272 the United States Food and Drug Administration which contains
273 Salvinorin A or its isomers, esters, ethers, salts, and salts of
274 isomers, esters, and ethers, if the existence of such isomers,
275 esters, ethers, and salts is possible within the specific
276 chemical designation.

277 37. Xylazine, except for a xylazine animal drug product
278 approved by the United States Food and Drug Administration and
279 the use of which conforms to the approved application or is
280 authorized under 21 U.S.C. s. 360b(a)(4). The manufacture,
281 importation, distribution, prescribing, or sale of xylazine for
282 human use is not subject to this exception.

283 38. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine)
284 (Thiophene analog of phencyclidine).

285 39. 3,4,5-Trimethoxyamphetamine.

286 40. Methyldone (3,4-Methylenedioxymethcathinone).

287 41. MDPV (3,4-Methylenedioxypyrovalerone).

288 42. Methyldmethcathinone.

289 43. Methoxymethcathinone.

290 44. Fluoromethcathinone.

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- 291 45. Methylethcathinone.
- 292 46. CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-
293 yl)phenol) and its dimethyloctyl (C8) homologue.
- 294 47. HU-210 [(6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-
295 methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol].
- 296 48. JWH-018 (1-Pentyl-3-(1-naphthoyl)indole).
- 297 49. JWH-073 (1-Butyl-3-(1-naphthoyl)indole).
- 298 50. JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-
299 naphthoyl)indole).
- 300 51. BZP (Benzylpiperazine).
- 301 52. Fluorophenylpiperazine.
- 302 53. Methylphenylpiperazine.
- 303 54. Chlorophenylpiperazine.
- 304 55. Methoxyphenylpiperazine.
- 305 56. DBZP (1,4-Dibenzylpiperazine).
- 306 57. TFMPP (Trifluoromethylphenylpiperazine).
- 307 58. MBDB (Methylbenzodioxolylbutanamine) or (3,4-
308 Methylenedioxy-N-methylbutanamine).
- 309 59. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
- 310 60. 5-Hydroxy-N-methyltryptamine.
- 311 61. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).
- 312 62. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
- 313 63. Methyltryptamine.
- 314 64. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
- 315 65. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
- 316 66. Tyramine (4-Hydroxyphenethylamine).
- 317 67. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
- 318 68. DiPT (N,N-Diisopropyltryptamine).
- 319 69. DPT (N,N-Dipropyltryptamine).

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- 320 70. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
 321 71. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
 322 72. DOI (4-Iodo-2,5-dimethoxyamphetamine).
 323 73. DOC (4-Chloro-2,5-dimethoxyamphetamine).
 324 74. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
 325 75. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
 326 76. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
 327 77. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
 328 78. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
 329 79. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
 330 80. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
 331 81. Butylone (3,4-Methylenedioxy-alpha-
 332 methylaminobutyrophenone).
 333 82. Ethcathinone.
 334 83. Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
 335 84. Naphyrone (Naphthylpyrovalerone).
 336 85. Dimethylone (3,4-Methylenedioxy-N,N-dimethylcathinone).
 337 86. 3,4-Methylenedioxy-N,N-diethylcathinone.
 338 87. 3,4-Methylenedioxy-propiofenone.
 339 88. 3,4-Methylenedioxy-alpha-bromopropiofenone.
 340 89. 3,4-Methylenedioxy-propiofenone-2-oxime.
 341 90. 3,4-Methylenedioxy-N-acetylcathinone.
 342 91. 3,4-Methylenedioxy-N-acetylmethcathinone.
 343 92. 3,4-Methylenedioxy-N-acetylethcathinone.
 344 93. Bromomethcathinone.
 345 94. Buphedrone (alpha-Methylamino-butyrophenone).
 346 95. Eutylone (3,4-Methylenedioxy-alpha-
 347 ethylaminobutyrophenone).
 348 96. Dimethylcathinone.

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97. Dimethylmethcathinone.

98. Pentylone (3,4-Methylenedioxy-alpha-methylaminovalerophenone).

99. MDPPP (3,4-Methylenedioxy-alpha-pyrrolidinopropiophenone).

100. MDPBP (3,4-Methylenedioxy-alpha-pyrrolidinobutyrophenone).

101. MOPPP (Methoxy-alpha-pyrrolidinopropiophenone).

102. MPHP (Methyl-alpha-pyrrolidinohexanophenone).

103. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP (Benocyclidine).

104. F-MABP (Fluoromethylaminobutyrophenone).

105. MeO-PBP (Methoxypyrrolidinobutyrophenone).

106. Et-PBP (Ethylpyrrolidinobutyrophenone).

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

108. Me-EABP (Methylethylaminobutyrophenone).

109. Etizolam.

110. PPP (Pyrrolidinopropiophenone).

111. PBP (Pyrrolidinobutyrophenone).

112. PVP (Pyrrolidinovalerophenone) or (Pyrrolidinopentiophenone).

113. MPPP (Methyl-alpha-pyrrolidinopropiophenone).

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

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

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

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

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

119. JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl)indole).

120. JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).

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

122. JWH-175 (1-Pentyl-3-(1-naphthylmethyl)indole).

123. JWH-201 (1-Pentyl-3-(4-methoxyphenylacetyl)indole).

124. JWH-203 (1-Pentyl-3-(2-chlorophenylacetyl)indole).

125. JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).

126. JWH-250 (1-Pentyl-3-(2-methoxyphenylacetyl)indole).

127. JWH-251 (1-Pentyl-3-(2-methylphenylacetyl)indole).

128. JWH-302 (1-Pentyl-3-(3-methoxyphenylacetyl)indole).

129. JWH-398 (1-Pentyl-3-(4-chloro-1-naphthoyl)indole).

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

131. HU-308 ([(1R,2R,5R)-2-[2,6-Dimethoxy-4-(2-methyloctan-2-yl)phenyl]-7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl]methanol).

132. HU-331 (3-Hydroxy-2-[(1R,6R)-3-methyl-6-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-2,5-cyclohexadiene-1,4-dione).

133. CB-13 (4-Pentyloxy-1-(1-naphthoyl)naphthalene).

134. CB-25 (N-Cyclopropyl-11-(3-hydroxy-5-pentylphenoxy)-undecanamide).

135. CB-52 (N-Cyclopropyl-11-(2-hexyl-5-hydroxyphenoxy)-undecanamide).

136. CP 55,940 (2-[3-Hydroxy-6-propanol-cyclohexyl]-5-(2-methyloctan-2-yl)phenol).

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

138. AM-2201 (1-(5-Fluoropentyl)-3-(1-naphthoyl)indole).

139. RCS-4 (1-Pentyl-3-(4-methoxybenzoyl)indole).

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140. RCS-8 (1-(2-Cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole).

141. WIN55,212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-naphthalenylmethanone).

142. WIN55,212-3 ([(3S)-2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-naphthalenylmethanone).

143. Pentedrone (alpha-Methylaminovalerophenone).

144. Fluoroamphetamine.

145. Fluoromethamphetamine.

146. Methoxetamine.

147. Methiopropamine.

148. Methylbuphedrone (Methyl-alpha-methylaminobutyrophenone).

149. APB ((2-Aminopropyl)benzofuran).

150. APDB ((2-Aminopropyl)-2,3-dihydrobenzofuran).

151. UR-144 (1-Pentyl-3-(2,2,3,3-tetramethylcyclopropanoyl)indole).

152. XLR11 (1-(5-Fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropanoyl)indole).

153. Chloro UR-144 (1-(Chloropentyl)-3-(2,2,3,3-tetramethylcyclopropanoyl)indole).

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

155. AM-2233(1-[(N-Methyl-2-piperidiny)methyl]-3-(2-iodobenzoyl)indole).

156. STS-135 (N-Adamant-1-yl 1-(5-fluoropentyl)indole-3-carboxamide).

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

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cyclohexylcarbamate).

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

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

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

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

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

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

164. 25I-NBOMe (4-Iodo-2,5-dimethoxy-[N-(2-
methoxybenzyl)]phenethylamine).

165. MDMA (3,4-Methylenedioxymethamphetamine).

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

167. Fluoro PB-22 (8-Quinoliny 1-(fluoropentyl)indole-3-
carboxylate).

168. BB-22 (8-Quinoliny 1-(cyclohexylmethyl)indole-3-
carboxylate).

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

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

171. AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-
(4-fluorobenzyl)indazole-3-carboxamide).

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

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

174. 25B-NBOMe (4-Bromo-2,5-dimethoxy-[N-(2-
methoxybenzyl)]phenethylamine).

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175. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

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

177. FUB-PB-22 (8-Quinoliny 1-(4-fluorobenzyl)indole-3-carboxylate).

178. Fluoro-NNEI (N-Naphthalen-1-yl 1-(fluoropentyl)indole-3-carboxamide).

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

180. THJ-2201 (1-(5-Fluoropentyl)-3-(1-naphthoyl)indazole).

181. AM-855 ((4aR,12bR)-8-Hexyl-2,5,5-trimethyl-1,4,4a,8,9,10,11,12b-octahydronaphtho[3,2-c]isochromen-12-ol).

182. AM-905 ((6aR,9R,10aR)-3-[(E)-Hept-1-enyl]-9-(hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-hexahydrobenzo[c]chromen-1-ol).

183. AM-906 ((6aR,9R,10aR)-3-[(Z)-Hept-1-enyl]-9-(hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-hexahydrobenzo[c]chromen-1-ol).

184. AM-2389 ((6aR,9R,10aR)-3-(1-Hexyl-cyclobut-1-yl)-6a,7,8,9,10,10a-hexahydro-6,6-dimethyl-6H-dibenzo[b,d]pyran-1,9-diol).

185. HU-243 ((6aR,8S,9S,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-8,9-ditritio-7,8,10,10a-tetrahydro-6aH-benzo[c]chromen-1-ol).

186. HU-336 ((6aR,10aR)-6,6,9-Trimethyl-3-pentyl-6a,7,10,10a-tetrahydro-1H-benzo[c]chromene-1,4(6H)-dione).

187. MAPB ((2-Methylaminopropyl)benzofuran).

188. 5-IT (2-(1H-Indol-5-yl)-1-methyl-ethylamine).

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494 189. 6-IT (2-(1H-Indol-6-yl)-1-methyl-ethylamine).

495 190. Synthetic Cannabinoids.—Unless specifically excepted
496 or unless listed in another schedule or contained within a
497 pharmaceutical product approved by the United States Food and
498 Drug Administration, any material, compound, mixture, or
499 preparation that contains any quantity of a synthetic
500 cannabinoid found to be in any of the following chemical class
501 descriptions, or homologues, nitrogen-heterocyclic analogs,
502 isomers (including optical, positional, or geometric), esters,
503 ethers, salts, and salts of homologues, nitrogen-heterocyclic
504 analogs, isomers, esters, or ethers, whenever the existence of
505 such homologues, nitrogen-heterocyclic analogs, isomers, esters,
506 ethers, salts, and salts of isomers, esters, or ethers is
507 possible within the specific chemical class or designation.
508 Since nomenclature of these synthetically produced cannabinoids
509 is not internationally standardized and may continually evolve,
510 these structures or the compounds of these structures shall be
511 included under this subparagraph, regardless of their specific
512 numerical designation of atomic positions covered, if it can be
513 determined through a recognized method of scientific testing or
514 analysis that the substance contains properties that fit within
515 one or more of the following categories:

516 a. Tetrahydrocannabinols.—Any tetrahydrocannabinols
517 naturally contained in a plant of the genus *Cannabis*, the
518 synthetic equivalents of the substances contained in the plant
519 or in the resinous extracts of the genus *Cannabis*, or synthetic
520 substances, derivatives, and their isomers with similar chemical
521 structure and pharmacological activity, including, but not
522 limited to, Delta 9 tetrahydrocannabinols and their optical

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

(I) Tetrahydrocannabinol.

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

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

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

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

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

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

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

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

(X) Parahexyl.

b. Naphthoylindoles, Naphthoylindazoles, Naphthoylcarbazoles, Naphthylmethylindoles, Naphthylmethylindazoles, and Naphthylmethylcarbazoles.—Any

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compound containing a naphthoylindole, naphthoylindazole, naphthoylcarbazole, naphthylmethylindole, naphthylmethylindazole, or naphthylmethylcarbazole structure, with or without substitution on the indole, indazole, or carbazole ring to any extent, whether or not substituted on the naphthyl ring to any extent, including, but not limited to:

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

(II) JWH-011 (1-(1-Methylhexyl)-2-methyl-3-(1-naphthoyl)indole).

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

(IV) JWH-016 (1-Butyl-2-methyl-3-(1-naphthoyl)indole).

(V) JWH-018 (1-Pentyl-3-(1-naphthoyl)indole).

(VI) JWH-019 (1-Hexyl-3-(1-naphthoyl)indole).

(VII) JWH-020 (1-Heptyl-3-(1-naphthoyl)indole).

(VIII) JWH-022 (1-(4-Pentenyl)-3-(1-naphthoyl)indole).

(IX) JWH-071 (1-Ethyl-3-(1-naphthoyl)indole).

(X) JWH-072 (1-Propyl-3-(1-naphthoyl)indole).

(XI) JWH-073 (1-Butyl-3-(1-naphthoyl)indole).

(XII) JWH-080 (1-Butyl-3-(4-methoxy-1-naphthoyl)indole).

(XIII) JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl)indole).

(XIV) JWH-098 (1-Pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole).

(XV) JWH-116 (1-Pentyl-2-ethyl-3-(1-naphthoyl)indole).

(XVI) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).

(XVII) JWH-149 (1-Pentyl-2-methyl-3-(4-methyl-1-naphthoyl)indole).

(XVIII) JWH-164 (1-Pentyl-3-(7-methoxy-1-naphthoyl)indole).

(XIX) JWH-175 (1-Pentyl-3-(1-naphthylmethyl)indole).

(XX) JWH-180 (1-Propyl-3-(4-propyl-1-naphthoyl)indole).

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581 (XXI) JWH-182 (1-Pentyl-3-(4-propyl-1-naphthoyl)indole).
582 (XXII) JWH-184 (1-Pentyl-3-[(4-methyl)-1-
583 naphthylmethyl]indole).
584 (XXIII) JWH-193 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methyl-1-
585 naphthoyl)indole).
586 (XXIV) JWH-198 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methoxy-1-
587 naphthoyl)indole).
588 (XXV) JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-
589 naphthoyl)indole).
590 (XXVI) JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).
591 (XXVII) JWH-387 (1-Pentyl-3-(4-bromo-1-naphthoyl)indole).
592 (XXVIII) JWH-398 (1-Pentyl-3-(4-chloro-1-naphthoyl)indole).
593 (XXIX) JWH-412 (1-Pentyl-3-(4-fluoro-1-naphthoyl)indole).
594 (XXX) JWH-424 (1-Pentyl-3-(8-bromo-1-naphthoyl)indole).
595 (XXXI) AM-1220 (1-[(1-Methyl-2-piperidinyl)methyl]-3-(1-
596 naphthoyl)indole).
597 (XXXII) AM-1235 (1-(5-Fluoropentyl)-6-nitro-3-(1-
598 naphthoyl)indole).
599 (XXXIII) AM-2201 (1-(5-Fluoropentyl)-3-(1-
600 naphthoyl)indole).
601 (XXXIV) Chloro JWH-018 (1-(Chloropentyl)-3-(1-
602 naphthoyl)indole).
603 (XXXV) Bromo JWH-018 (1-(Bromopentyl)-3-(1-
604 naphthoyl)indole).
605 (XXXVI) AM-2232 (1-(4-Cyanobutyl)-3-(1-naphthoyl)indole).
606 (XXXVII) THJ-2201 (1-(5-Fluoropentyl)-3-(1-
607 naphthoyl)indazole).
608 (XXXVIII) MAM-2201 (1-(5-Fluoropentyl)-3-(4-methyl-1-
609 naphthoyl)indole).

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

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

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

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

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

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

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

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

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

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

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

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

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

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

d. Naphthylmethylenindenes.—Any compound containing a naphthylmethylenindene structure, with or without substitution at the 3-position of the indene ring to any extent, whether or not substituted on the naphthyl ring to any extent, including, but not limited to, JWH-176 (3-Pentyl-1-

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(naphthylmethylen)indene).

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

(I) JWH-167 (1-Pentyl-3-(phenylacetyl)indole).

(II) JWH-201 (1-Pentyl-3-(4-methoxyphenylacetyl)indole).

(III) JWH-203 (1-Pentyl-3-(2-chlorophenylacetyl)indole).

(IV) JWH-250 (1-Pentyl-3-(2-methoxyphenylacetyl)indole).

(V) JWH-251 (1-Pentyl-3-(2-methylphenylacetyl)indole).

(VI) JWH-302 (1-Pentyl-3-(3-methoxyphenylacetyl)indole).

(VII) Cannabipiperidiethanone.

(VIII) RCS-8 (1-(2-Cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole).

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

(I) CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-yl)phenol).

(II) Cannabicyclohexanol (CP 47,497 dimethyloctyl (C8) homologue).

(III) CP-55,940 (2-(3-Hydroxy-6-propanol-cyclohexyl)-5-(2-methyloctan-2-yl)phenol).

g. Benzoylindoles and Benzoylindazoles.—Any compound containing a benzoylindole or benzoylindazole structure, with or without substitution on the indole or indazole ring to any

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extent, whether or not substituted on the phenyl ring to any extent, including, but not limited to:

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

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

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

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

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

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

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

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

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

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

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

(III) Chloro UR-144 (1-(Chloropentyl)-3-(2,2,3,3-tetramethylcyclopropanoyl)indole).

(IV) A-796,260 (1-[2-(4-Morpholinyl)ethyl]-3-(2,2,3,3-

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tetramethylcyclopropanoyl)indole).

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

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

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

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

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

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

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

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

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

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

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

(V) AB-001 (1-Pentyl-3-(1-adamantoyl)indole).

(VI) APICA (N-Adamant-1-yl 1-pentylindole-3-carboxamide).

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(VII) Fluoro AB-001 (1-(Fluoropentyl)-3-(1-adamantoyl)indole).

j. Quinolinyndolecarboxylates, Quinolinyndazolecarboxylates, Quinolinyndazolecarboxamides, and Quinolinyndazolecarboxamides.—Any compound containing a quinolinyndole carboxylate, quinolinyndazole carboxylate, isoquinolinyndole carboxylate, isoquinolinyndazole carboxylate, quinolinyndole carboxamide, quinolinyndazole carboxamide, isoquinolinyndole carboxamide, or isoquinolinyndazole carboxamide structure, with or without substitution on the indole or indazole ring to any extent, whether or not substituted on the quinoline or isoquinoline ring to any extent, including, but not limited to:

(I) PB-22 (8-Quinolinyndyl 1-pentyndazole-3-carboxylate).

(II) Fluoro PB-22 (8-Quinolinyndyl 1-(fluoropentyl)indole-3-carboxylate).

(III) BB-22 (8-Quinolinyndyl 1-(cyclohexylmethyl)indole-3-carboxylate).

(IV) FUB-PB-22 (8-Quinolinyndyl 1-(4-fluorobenzyl)indole-3-carboxylate).

(V) NPB-22 (8-Quinolinyndyl 1-pentyndazole-3-carboxylate).

(VI) Fluoro NPB-22 (8-Quinolinyndyl 1-(fluoropentyl)indazole-3-carboxylate).

(VII) FUB-NPB-22 (8-Quinolinyndyl 1-(4-fluorobenzyl)indazole-3-carboxylate).

(VIII) THJ (8-Quinolinyndyl 1-pentyndazole-3-carboxamide).

(IX) Fluoro THJ (8-Quinolinyndyl 1-(fluoropentyl)indazole-3-carboxamide).

k. Naphthylndolecarboxylates and

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Naphthylindazolecarboxylates.—Any compound containing a naphthylindole carboxylate or naphthylindazole carboxylate structure, with or without substitution on the indole or indazole ring to any extent, whether or not substituted on the naphthyl ring to any extent, including, but not limited to:

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

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

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

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

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

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

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

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

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

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

(V) Fluoro MN-18 (N-Naphthalen-1-yl 1-

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(fluoropentyl)indazole-3-carboxamide).

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

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

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

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

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

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

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

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

(VIII) AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).

(IX) ADB-FUBINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-

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yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

191. Substituted Cathinones.—Unless specifically excepted,
listed in another schedule, or contained within a pharmaceutical
product approved by the United States Food and Drug

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Administration, any material, compound, mixture, or preparation, including its salts, isomers, esters, or ethers, and salts of isomers, esters, or ethers, whenever the existence of such salts is possible within any of the following specific chemical designations:

- a. Any compound containing a 2-amino-1-phenyl-1-propanone structure;
- b. Any compound containing a 2-amino-1-naphthyl-1-propanone structure; or
- c. Any compound containing a 2-amino-1-thiophenyl-1-propanone structure,

whether or not the compound is further modified:

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

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

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

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

- (A) Methcathinone.
- (B) Ethcathinone.
- (C) Methylone (3,4-Methylenedioxymethcathinone).

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- 900 (D) 2,3-Methylenedioxymethcathinone.
- 901 (E) MDPV (3,4-Methylenedioxypyrovalerone).
- 902 (F) Methylnmethcathinone.
- 903 (G) Methoxymethcathinone.
- 904 (H) Fluoromethcathinone.
- 905 (I) Methylethcathinone.
- 906 (J) Butylone (3,4-Methylenedioxy-alpha-
- 907 methylaminobutyrophenone).
- 908 (K) Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
- 909 (L) BMDP (3,4-Methylenedioxy-N-benzylcathinone).
- 910 (M) Naphyrone (Naphthylpyrovalerone).
- 911 (N) Bromomethcathinone.
- 912 (O) Buphedrone (alpha-Methylaminobutyrophenone).
- 913 (P) Eutylone (3,4-Methylenedioxy-alpha-
- 914 ethylaminobutyrophenone).
- 915 (Q) Dimethylcathinone.
- 916 (R) Dimethylmethcathinone.
- 917 (S) Pentylone (3,4-Methylenedioxy-alpha-
- 918 methylaminovalerophenone).
- 919 (T) Pentedrone (alpha-Methylaminovalerophenone).
- 920 (U) MDPPP (3,4-Methylenedioxy-alpha-
- 921 pyrrolidinopropiophenone).
- 922 (V) MDPBP (3,4-Methylenedioxy-alpha-
- 923 pyrrolidinobutyrophenone).
- 924 (W) MPPP (Methyl-alpha-pyrrolidinopropiophenone).
- 925 (X) PPP (Pyrrolidinopropiophenone).
- 926 (Y) PVP (Pyrrolidinovalerophenone) or
- 927 (Pyrrolidinopentiophenone).
- 928 (Z) MOPPP (Methoxy-alpha-pyrrolidinopropiophenone).

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(AA) MPHP (Methyl-alpha-pyrrolidinohexanophenone).
 (BB) F-MABP (Fluoromethylaminobutyrophenone).
 (CC) Me-EABP (Methylethylaminobutyrophenone).
 (DD) PBP (Pyrrolidinobutyrophenone).
 (EE) MeO-PBP (Methoxypyrrolidinobutyrophenone).
 (FF) Et-PBP (Ethylpyrrolidinobutyrophenone).
 (GG) 3-Me-4-MeO-MCAT (3-Methyl-4-Methoxymethcathinone).
 (HH) Dimethylone (3,4-Methylenedioxy-N,N-dimethylcathinone).
 (II) 3,4-Methylenedioxy-N,N-diethylcathinone.
 (JJ) 3,4-Methylenedioxy-N-acetylcathinone.
 (KK) 3,4-Methylenedioxy-N-acetylmethcathinone.
 (LL) 3,4-Methylenedioxy-N-acetylethcathinone.
 (MM) Methylbuphedrone (Methyl-alpha-methylaminobutyrophenone).
 (NN) Methyl-alpha-methylaminohexanophenone.
 (OO) N-Ethyl-N-methylcathinone.
 (PP) PHP (Pyrrolidinohexanophenone).
 (QQ) PV8 (Pyrrolidinoheptanophenone).
 (RR) Chloromethcathinone.
 (SS) 4-Bromo-2,5-dimethoxy-alpha-aminoacetophenone.
 192. Substituted Phenethylamines.—Unless specifically
 excepted or unless listed in another schedule, or contained
 within a pharmaceutical product approved by the United States
 Food and Drug Administration, any material, compound, mixture,
 or preparation, including its salts, isomers, esters, or ethers,
 and salts of isomers, esters, or ethers, whenever the existence
 of such salts is possible within any of the following specific
 chemical designations, any compound containing a phenethylamine

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structure, without a beta-keto group, and without a benzyl group attached to the amine group, whether or not the compound is further modified with or without substitution on the phenyl ring to any extent with alkyl, alkylthio, nitro, alkoxy, thio, halide, fused alkylenedioxy, fused furan, fused benzofuran, fused dihydrofuran, or fused tetrahydropyran substituents, whether or not further substituted on a ring to any extent, with or without substitution at the alpha or beta position by any alkyl substituent, with or without substitution at the nitrogen atom, and with or without inclusion of the 2-amino nitrogen atom in a cyclic structure, including, but not limited to:

- a. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
- b. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
- c. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
- d. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
- e. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
- f. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
- g. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
- h. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
- i. 2C-D (4-Methyl-2,5-dimethoxyphenethylamine).
- j. 2C-H (2,5-Dimethoxyphenethylamine).
- k. 2C-N (4-Nitro-2,5-dimethoxyphenethylamine).
- l. 2C-P (4-(n)-Propyl-2,5-dimethoxyphenethylamine).
- m. MDMA (3,4-Methylenedioxyamphetamine).
- n. MBDB (Methylbenzodioxolylbutanamine) or (3,4-Methylenedioxy-N-methylbutanamine).
- o. MDA (3,4-Methylenedioxyamphetamine).
- p. 2,5-Dimethoxyamphetamine.
- q. Fluoroamphetamine.

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987 r. Fluoromethamphetamine.
 988 s. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
 989 t. DOB (4-Bromo-2,5-dimethoxyamphetamine).
 990 u. DOC (4-Chloro-2,5-dimethoxyamphetamine).
 991 v. DOET (4-Ethyl-2,5-dimethoxyamphetamine).
 992 w. DOI (4-Iodo-2,5-dimethoxyamphetamine).
 993 x. DOM (4-Methyl-2,5-dimethoxyamphetamine).
 994 y. PMA (4-Methoxyamphetamine).
 995 z. N-Ethylamphetamine.
 996 aa. 3,4-Methylenedioxy-N-hydroxyamphetamine.
 997 bb. 5-Methoxy-3,4-methylenedioxyamphetamine.
 998 cc. PMMA (4-Methoxymethamphetamine).
 999 dd. N,N-Dimethylamphetamine.
 1000 ee. 3,4,5-Trimethoxyamphetamine.
 1001 ff. 4-APB (4-(2-Aminopropyl)benzofuran).
 1002 gg. 5-APB (5-(2-Aminopropyl)benzofuran).
 1003 hh. 6-APB (6-(2-Aminopropyl)benzofuran).
 1004 ii. 7-APB (7-(2-Aminopropyl)benzofuran).
 1005 jj. 4-APDB (4-(2-Aminopropyl)-2,3-dihydrobenzofuran).
 1006 kk. 5-APDB (5-(2-Aminopropyl)-2,3-dihydrobenzofuran).
 1007 ll. 6-APDB (6-(2-Aminopropyl)-2,3-dihydrobenzofuran).
 1008 mm. 7-APDB (7-(2-Aminopropyl)-2,3-dihydrobenzofuran).
 1009 nn. 4-MAPB (4-(2-Methylaminopropyl)benzofuran).
 1010 oo. 5-MAPB (5-(2-Methylaminopropyl)benzofuran).
 1011 pp. 6-MAPB (6-(2-Methylaminopropyl)benzofuran).
 1012 qq. 7-MAPB (7-(2-Methylaminopropyl)benzofuran).
 1013 rr. 5-EAPB (5-(2-Ethylaminopropyl)benzofuran).
 1014 ss. 5-MAPDB (5-(2-Methylaminopropyl)-2,3-
 1015 dihydrobenzofuran),

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

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

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

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

c. 25B-NBF (4-Bromo-2,5-dimethoxy-[N-(2-fluorobenzyl)]phenethylamine).

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d. 25B-NBMD (4-Bromo-2,5-dimethoxy-[N-(2,3-methylenedioxybenzyl)]phenethylamine).

e. 25I-NBOMe (4-Iodo-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

f. 25I-NBOH (4-Iodo-2,5-dimethoxy-[N-(2-hydroxybenzyl)]phenethylamine).

g. 25I-NBF (4-Iodo-2,5-dimethoxy-[N-(2-fluorobenzyl)]phenethylamine).

h. 25I-NBMD (4-Iodo-2,5-dimethoxy-[N-(2,3-methylenedioxybenzyl)]phenethylamine).

i. 25T2-NBOMe (4-Methylthio-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

j. 25T4-NBOMe (4-Isopropylthio-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

k. 25T7-NBOMe (4-(n)-Propylthio-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

l. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

m. 25C-NBOH (4-Chloro-2,5-dimethoxy-[N-(2-hydroxybenzyl)]phenethylamine).

n. 25C-NBF (4-Chloro-2,5-dimethoxy-[N-(2-fluorobenzyl)]phenethylamine).

o. 25C-NBMD (4-Chloro-2,5-dimethoxy-[N-(2,3-methylenedioxybenzyl)]phenethylamine).

p. 25H-NBOMe (2,5-Dimethoxy-[N-(2-methoxybenzyl)]phenethylamine).

q. 25H-NBOH (2,5-Dimethoxy-[N-(2-hydroxybenzyl)]phenethylamine).

r. 25H-NBF (2,5-Dimethoxy-[N-(2-

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fluorobenzyl)]phenethylamine).

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

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

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

- a. Alpha-Ethyltryptamine.
- b. Bufotenine.
- c. DET (Diethyltryptamine).
- d. DMT (Dimethyltryptamine).
- e. MET (N-Methyl-N-ethyltryptamine).
- f. DALT (N,N-Diallyltryptamine).
- g. EiPT (N-Ethyl-N-isopropyltryptamine).
- h. MiPT (N-Methyl-N-isopropyltryptamine).
- i. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
- j. 5-Hydroxy-N-methyltryptamine.
- k. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).

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1. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
 m. Methyltryptamine.
 n. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
 o. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
 p. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
 q. DiPT (N,N-Diisopropyltryptamine).
 r. DPT (N,N-Dipropyltryptamine).
 s. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
 t. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
 u. 4-AcO-DMT (4-Acetoxy-N,N-dimethyltryptamine).
 v. 4-AcO-DiPT (4-Acetoxy-N,N-diisopropyltryptamine).
 w. 4-Hydroxy-DET (4-Hydroxy-N,N-diethyltryptamine).
 x. 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-ethyltryptamine).
 y. 4-Hydroxy-MiPT (4-Hydroxy-N-methyl-N-
 isopropyltryptamine).
 z. Methyl-alpha-ethyltryptamine.
 aa. Bromo-DALT (Bromo-N,N-diallyltryptamine),
 which does not include tryptamine, psilocyn as described in
 subparagraph 34., or psilocybin as described in subparagraph 33.
 195. Substituted Phenylcyclohexylamines.—Unless
 specifically excepted or unless listed in another schedule, or
 contained within a pharmaceutical product approved by the United
 States Food and Drug Administration, any material, compound,
 mixture, or preparation containing a phenylcyclohexylamine
 structure, with or without any substitution on the phenyl ring,
 any substitution on the cyclohexyl ring, any replacement of the
 phenyl ring with a thiophenyl or benzothiophenyl ring, with or
 without substitution on the amine with alkyl, dialkyl, or alkoxy

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substituents, inclusion of the nitrogen in a cyclic structure,
or any combination of the above, including, but not limited to:

- a. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP (Benocyclidine).
- b. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine analog of phencyclidine).
- c. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine analog of phencyclidine).
- d. PCPr (Phenylcyclohexylpropylamine).
- e. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine) (Thiophene analog of phencyclidine).
- f. PCEEA (Phenylcyclohexyl(ethoxyethylamine)).
- g. PCMPA (Phenylcyclohexyl(methoxypropylamine)).
- h. Methoxetamine.
- i. 3-Methoxy-PCE ((3-Methoxyphenyl)cyclohexylethylamine).
- j. Bromo-PCP ((Bromophenyl)cyclohexylpiperidine).
- k. Chloro-PCP ((Chlorophenyl)cyclohexylpiperidine).
- l. Fluoro-PCP ((Fluorophenyl)cyclohexylpiperidine).
- m. Hydroxy-PCP ((Hydroxyphenyl)cyclohexylpiperidine).
- n. Methoxy-PCP ((Methoxyphenyl)cyclohexylpiperidine).
- o. Methyl-PCP ((Methylphenyl)cyclohexylpiperidine).
- p. Nitro-PCP ((Nitrophenyl)cyclohexylpiperidine).
- q. Oxo-PCP ((Oxophenyl)cyclohexylpiperidine).
- r. Amino-PCP ((Aminophenyl)cyclohexylpiperidine).
196. W-15, 4-chloro-N-[1-(2-phenylethyl)-2-piperidinylidene]-benzenesulfonamide.
197. W-18, 4-chloro-N-[1-[2-(4-nitrophenyl)ethyl]-2-piperidinylidene]-benzenesulfonamide.
198. AH-7921, 3,4-dichloro-N-[[1-

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(dimethylamino)cyclohexyl)methyl]-benzamide.

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

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

Section 2. Paragraph (i) of subsection (1) of section 893.13, Florida Statutes, is amended to read:

893.13 Prohibited acts; penalties.—

(1)

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

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

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

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

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

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

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

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

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

~~h.g.~~ A mixture containing any substance described in sub-subparagraphs a.-g. ~~a.-f.~~; and

2. The substance or mixture listed in subparagraph 1. is in

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a form that resembles, or is mixed, granulated, absorbed, spray-dried, or aerosolized as or onto, coated on, in whole or in part, or solubilized with or into, a product, when such product or its packaging further has at least one of the following attributes:

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

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

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

d. Contains a cartoon character imprint.

Section 3. Paragraph (c) of subsection (1) of section 893.135, Florida Statutes, is amended to read:

893.135 Trafficking; mandatory sentences; suspension or reduction of sentences; conspiracy to engage in trafficking.—

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

(c)1. A person who knowingly sells, purchases, manufactures, delivers, or brings into this state, or who is knowingly in actual or constructive possession of, 4 grams or more of any morphine, opium, hydromorphone, or any salt, derivative, isomer, or salt of an isomer thereof, including heroin, as described in s. 893.03(1)(b), (2)(a), (3)(c)3., or (3)(c)4., or 4 grams or more of any mixture containing any such substance, but less than 30 kilograms of such substance or mixture, commits a felony of the first degree, which felony shall be known as "trafficking in illegal drugs," punishable as provided in s. 775.082, s. 775.083, or s. 775.084. If the

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quantity involved:

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

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

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

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

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

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

c. Is 100 grams or more, but less than 300 grams, such person shall be sentenced to a mandatory minimum term of

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imprisonment of 15 years and shall be ordered to pay a fine of \$500,000.

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

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

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

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

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

d. Is 100 grams or more, but less than 30 kilograms, such person shall be sentenced to a mandatory minimum term of imprisonment of 25 years and shall be ordered to pay a fine of \$750,000.

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1277 4.a. A person who knowingly sells, purchases, manufactures,
1278 delivers, or brings into this state, or who is knowingly in
1279 actual or constructive possession of, 4 grams or more of:

1280 (I) Alfentanil, as described in s. 893.03(2)(b)1.;

1281 (II) Carfentanil, as described in s. 893.03(2)(b)6.;

1282 (III) Fentanyl, as described in s. 893.03(2)(b)9.;

1283 (IV) Sufentanil, as described in s. 893.03(2)(b)30.;

1284 (V) A fentanyl derivative, as described in s.

1285 893.03(1)(a)63.;

1286 (VI) A controlled substance analog, as described in s.

1287 893.0356, of any substance described in sub-sub-subparagraphs

1288 (I)-(V); or

1289 (VII) A mixture containing any substance described in sub-
1290 sub-subparagraphs (I)-(VI),

1291
1292 commits a felony of the first degree, which felony shall be
1293 known as "trafficking in dangerous fentanyl or fentanyl
1294 analogues," punishable as provided in s. 775.082, s. 775.083, or
1295 s. 775.084.

1296 b. If the quantity involved under sub-subparagraph a.:

1297 (I) Is 4 grams or more, but less than 14 grams, such person
1298 shall be sentenced to a mandatory minimum term of imprisonment
1299 of 7 years~~7~~ and shall be ordered to pay a fine of \$50,000.

1300 (II) Is 14 grams or more, but less than 28 grams, such
1301 person shall be sentenced to a mandatory minimum term of
1302 imprisonment of 20 years~~7~~ and shall be ordered to pay a fine of
1303 \$100,000.

1304 (III) Is 28 grams or more, such person shall be sentenced
1305 to a mandatory minimum term of imprisonment of 25 years~~7~~ and

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shall be ordered to pay a fine of \$500,000.

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

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

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

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

(IV) Contains a cartoon character imprint.

5. A person who knowingly sells, purchases, manufactures, delivers, or brings into this state, or who is knowingly in actual or constructive possession of, 30 kilograms or more of any morphine, opium, oxycodone, hydrocodone, codeine, hydromorphone, or any salt, derivative, isomer, or salt of an isomer thereof, including heroin, as described in s. 893.03(1)(b), (2)(a), (3)(c)3., or (3)(c)4., or 30 kilograms or more of any mixture containing any such substance, commits the first degree felony of trafficking in illegal drugs. A person who has been convicted of the first degree felony of trafficking

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in illegal drugs under this subparagraph shall be punished by life imprisonment and is ineligible for any form of discretionary early release except pardon or executive clemency or conditional medical release under s. 947.149. However, if the court determines that, in addition to committing any act specified in this paragraph:

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

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

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

6. A person who knowingly brings into this state 60 kilograms or more of any morphine, opium, oxycodone, hydrocodone, codeine, hydromorphone, or any salt, derivative, isomer, or salt of an isomer thereof, including heroin, as described in s. 893.03(1)(b), (2)(a), (3)(c)3., or (3)(c)4., or 60 kilograms or more of any mixture containing any such substance, and who knows that the probable result of such importation would be the death of a person, commits capital importation of illegal drugs, a capital felony punishable as provided in ss. 775.082 and 921.142. A person sentenced for a capital felony under this paragraph shall also be sentenced to

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pay the maximum fine provided under subparagraph 1.

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

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

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

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

Section 4. Except as otherwise expressly provided in this act and except for this section, which shall take effect July 1, 2026, this act shall take effect October 1, 2026.