A bill to be entitled

An act relating to controlled substances; amending s. 893.03, F.S.; adding to the list of Schedule I controlled substances 7-Hydroxymitragynine concentrated at a level above 400 parts per million on a dry-weight basis; 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. 893.131, F.S.; conforming a cross-reference; 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 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.

9. Benzethidine.

10. Benzylfentanyl.

11. Betacetylmethadol.

12. Beta-hydroxyfentanyl.

13. Beta-hydroxy-3-methylfentanyl.

14. Betameprodine.

15. Betamethadol.

16. Betaprodine.

17. Clonitazene.

18. Dextromoramide.

19. Diampromide.

20. Diethylthiambutene.

21. Difenoxin.

22. Dimenoxadol.

23. Dimepheptanol.

24. Dimethylthiambutene.

25. Dioxaphetyl butyrate.

26. Dipipanone.

27. Ethylmethylthiambutene.

28. Etonitazene.

29. Etoxeridine.

30. Flunitrazepam.

31. Furethidine.

32. 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.

33.~~32.~~ Hydroxypethidine.

34.~~33.~~ Ketobemidone.

35.~~34.~~ Levomoramide.

36.~~35.~~ Levophenacylmorphan.

37.~~36.~~ Desmethylprodine (1-Methyl-4-Phenyl-4-Propionoxypiperidine).

38.~~37.~~ 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide).

39.~~38.~~ 3-Methylthiofentanyl.

40.~~39.~~ Morpheridine.

41.~~40.~~ Noracymethadol.

42.~~41.~~ Norlevorphanol.

43.~~42.~~ Normethadone.

44.~~43.~~ Norpipanone.

45.~~44.~~ Para-Fluorofentanyl.

46.~~45.~~ Phenadoxone.

47.~~46.~~ Phenampromide.

48.~~47.~~ Phenomorphan.

49.~~48.~~ Phenoperidine.

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

51.~~50.~~ Piritramide.

52.~~51.~~ Proheptazine.

53.~~52.~~ Properidine.

54.~~53.~~ Propiram.

55.~~54.~~ Racemoramide.

56.~~55.~~ Thenylfentanyl.

57.~~56.~~ Thiofentanyl.

58.~~57.~~ Tianeptine.

59.~~58.~~ Tilidine.

60.~~59.~~ Trimeperidine.

61.~~60.~~ Acetylfentanyl.

62.~~61.~~ Butyrylfentanyl.

63.~~62.~~ Beta-Hydroxythiofentanyl.

64.~~63.~~ Fentanyl 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 4-anilidopiperidine structure:

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

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

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

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

e. With or without substitution at the alpha or beta position of the piperidine ring with alkyl, hydroxyl, or methoxy

groups;

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

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.

(XXII) Isobutyryl fentanyl.

(XXIII) Remifentanil.

65.~~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.

(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.

(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).

5. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).

6. Bufotenine.

7. Cannabis.

8. Cathinone.

9. DET (Diethyltryptamine).

10. 2,5-Dimethoxyamphetamine.

11. DOET (4-Ethyl-2,5-Dimethoxyamphetamine).

12. DMT (Dimethyltryptamine).

13. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine analog of phencyclidine).

14. JB-318 (N-Ethyl-3-piperidyl benzilate).

15. N-Ethylamphetamine.

16. Fenethylline.

17. 3,4-Methylenedioxy-N-hydroxyamphetamine.

18. Ibogaine.

19. LSD (Lysergic acid diethylamide).

20. Mescaline.

21. Methcathinone.

22. 5-Methoxy-3,4-methylenedioxyamphetamine.

23. PMA (4-Methoxyamphetamine).

24. PMMA (4-Methoxymethamphetamine).

25. DOM (4-Methyl-2,5-dimethoxyamphetamine).

26. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).

27. MDA (3,4-Methylenedioxyamphetamine).

28. JB-336 (N-Methyl-3-piperidyl benzilate).

29. N,N-Dimethylamphetamine.

30. Parahexyl.

31. Peyote.

32. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine analog of phencyclidine).

33. Psilocybin.

34. Psilocyn.

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

36. Salvinorin A, except for any drug product approved by the United States Food and Drug Administration which contains Salvinorin A or its isomers, esters, ethers, salts, and salts of

isomers, esters, and ethers, if the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation.

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

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

39. 3,4,5-Trimethoxyamphetamine.

40. Methylone (3,4-Methylenedioxymethcathinone).

41. MDPV (3,4-Methylenedioxypyrovalerone).

42. Methylmethcathinone.

43. Methoxymethcathinone.

44. Fluoromethcathinone.

45. Methylethcathinone.

46. CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-yl)phenol) and its dimethyloctyl (C8) homologue.

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

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

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

50. JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-naphthoyl)indole).
51. BZP (Benzylpiperazine).
52. Fluorophenylpiperazine.
53. Methylphenylpiperazine.
54. Chlorophenylpiperazine.
55. Methoxyphenylpiperazine.
56. DBZP (1,4-Dibenzylpiperazine).
57. TFMPP (Trifluoromethylphenylpiperazine).
58. MBDB (Methylbenzodioxolylbutanamine) or (3,4-Methylenedioxy-N-methylbutanamine).
59. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
60. 5-Hydroxy-N-methyltryptamine.
61. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).
62. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
63. Methyltryptamine.
64. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
65. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
66. Tyramine (4-Hydroxyphenethylamine).
67. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
68. DiPT (N,N-Diisopropyltryptamine).
69. DPT (N,N-Dipropyltryptamine).
70. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
71. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
72. DOI (4-Iodo-2,5-dimethoxyamphetamine).

73. DOC (4-Chloro-2,5-dimethoxyamphetamine).
74. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
75. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
76. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
77. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
78. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
79. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
80. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
81. Butylone (3,4-Methylenedioxy-alpha-methylaminobutyrophenone).
82. Ethcathinone.
83. Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
84. Naphyrone (Naphthylpyrovalerone).
85. Dimethylone (3,4-Methylenedioxy-N,N-dimethylcathinone).
86. 3,4-Methylenedioxy-N,N-diethylcathinone.
87. 3,4-Methylenedioxy-propiophenone.
88. 3,4-Methylenedioxy-alpha-bromopropiophenone.
89. 3,4-Methylenedioxy-propiophenone-2-oxime.
90. 3,4-Methylenedioxy-N-acetylcathinone.
91. 3,4-Methylenedioxy-N-acetylmethcathinone.
92. 3,4-Methylenedioxy-N-acetylethcathinone.
93. Bromomethcathinone.
94. Buphedrone (alpha-Methylamino-butyrophenone).
95. Eutylone (3,4-Methylenedioxy-alpha-

ethylaminobutyrophenone).

96. Dimethylcathinone.

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).
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).

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-piperidinyl)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)-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-Quinolinyl 1-pentylindole-3-carboxylate).

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

168. BB-22 (8-Quinolinyl 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).

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-Quinolinyl 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).

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

190. Synthetic Cannabinoids.—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 that contains any quantity of a synthetic cannabinoid found to be in any of the following chemical class descriptions, or homologues, nitrogen-heterocyclic analogs, isomers (including optical, positional, or geometric), esters, ethers, salts, and salts of homologues, nitrogen-heterocyclic analogs, isomers, esters, or ethers, whenever the existence of such homologues, nitrogen-heterocyclic analogs, isomers, esters, ethers, salts, and salts of isomers, esters, or ethers is possible within the specific chemical class or designation. Since nomenclature of these synthetically produced cannabinoids is not internationally standardized and may continually evolve, these structures or the compounds of these structures shall be included under this subparagraph, regardless of their specific numerical designation of atomic positions covered, if it can be determined through a recognized method of scientific testing or analysis that the substance contains properties that fit within one or more of the following categories:

a. Tetrahydrocannabinols.—Any tetrahydrocannabinols naturally contained in a plant of the genus Cannabis, the synthetic equivalents of the substances contained in the plant or in the resinous extracts of the genus Cannabis, or synthetic

substances, derivatives, and their isomers with similar chemical structure and pharmacological activity, including, but not limited to, Delta 9 tetrahydrocannabinols and their optical 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 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).
(XXI) JWH-182 (1-Pentyl-3-(4-propyl-1-naphthoyl)indole).
(XXII) JWH-184 (1-Pentyl-3-[(4-methyl)-1-naphthylmethyl]indole).
(XXIII) JWH-193 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methyl-1-naphthoyl)indole).
(XXIV) JWH-198 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methoxy-1-naphthoyl)indole).
(XXV) JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-naphthoyl)indole).
(XXVI) JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).
(XXVII) JWH-387 (1-Pentyl-3-(4-bromo-1-naphthoyl)indole).
(XXVIII) JWH-398 (1-Pentyl-3-(4-chloro-1-

naphthoyl)indole).

(XXIX) JWH-412 (1-Pentyl-3-(4-fluoro-1-naphthoyl)indole).

(XXX) JWH-424 (1-Pentyl-3-(8-bromo-1-naphthoyl)indole).

(XXXI) AM-1220 (1-[(1-Methyl-2-piperidinyl)methyl]-3-(1-naphthoyl)indole).

(XXXII) AM-1235 (1-(5-Fluoropentyl)-6-nitro-3-(1-naphthoyl)indole).

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

(XXXIV) Chloro JWH-018 (1-(Chloropentyl)-3-(1-naphthoyl)indole).

(XXXV) Bromo JWH-018 (1-(Bromopentyl)-3-(1-naphthoyl)indole).

(XXXVI) AM-2232 (1-(4-Cyanobutyl)-3-(1-naphthoyl)indole).

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

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

(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-(naphthylmethylene)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 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-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).

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

j. Quinolinylindolecarboxylates, Quinolinylindazolecarboxylates, Quinolinylindolecarboxamides, and Quinolinylindazolecarboxamides.—Any compound containing a quinolinylindole carboxylate, quinolinylindazole carboxylate, isoquinolinylindole carboxylate, isoquinolinylindazole carboxylate, quinolinylindole carboxamide, quinolinylindazole carboxamide, isoquinolinylindole carboxamide, or isoquinolinylindazole 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-Quinolinyl 1-pentylindole-3-carboxylate).

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

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

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

(V) NPB-22 (8-Quinolinyl 1-pentylindazole-3-carboxylate).

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

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

(VIII) THJ (8-Quinolinyl 1-pentylindazole-3-carboxamide).

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

k. Naphthylindolecarboxylates and 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).

l. 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-(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-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).

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 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).

(D) 2,3-Methylenedioxymethcathinone.

(E) MDPV (3,4-Methylenedioxypyrovalerone).

(F) Methylmethcathinone.

(G) Methoxymethcathinone.

(H) Fluoromethcathinone.

(I) Methylethcathinone.

(J) Butylone (3,4-Methylenedioxy-alpha-methylaminobutyrophenone).

(K) Ethylone (3,4-Methylenedioxy-N-ethylcathinone).

(L) BMDP (3,4-Methylenedioxy-N-benzylcathinone).

(M) Naphyrone (Naphthylpyrovalerone).

(N) Bromomethcathinone.

(O) Buphedrone (alpha-Methylaminobutyrophenone).

(P) Eutylone (3,4-Methylenedioxy-alpha-ethylaminobutyrophenone).

(Q) Dimethylcathinone.

(R) Dimethylmethcathinone.

(S) Pentylone (3,4-Methylenedioxy-alpha-

methylaminovalerophenone).

(T) Pentedrone (alpha-Methylaminovalerophenone).

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

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

(W) MPPP (Methyl-alpha-pyrrolidinopropiophenone).

(X) PPP (Pyrrolidinopropiophenone).

(Y) PVP (Pyrrolidinovalerophenone) or (Pyrrolidinopentiophenone).

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

(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 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-Methylenedioxymethamphetamine).

n. MBDB (Methylbenzodioxolylbutanamine) or (3,4-Methylenedioxy-N-methylbutanamine).

o. MDA (3,4-Methylenedioxyamphetamine).

p. 2,5-Dimethoxyamphetamine.

q. Fluoroamphetamine.

r. Fluoromethamphetamine.

s. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).

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

u. DOC (4-Chloro-2,5-dimethoxyamphetamine).

v. DOET (4-Ethyl-2,5-dimethoxyamphetamine).

w. DOI (4-Iodo-2,5-dimethoxyamphetamine).

x. DOM (4-Methyl-2,5-dimethoxyamphetamine).

y. PMA (4-Methoxyamphetamine).

z. N-Ethylamphetamine.

aa. 3,4-Methylenedioxy-N-hydroxyamphetamine.

bb. 5-Methoxy-3,4-methylenedioxyamphetamine.

cc. PMMA (4-Methoxymethamphetamine).

dd. N,N-Dimethylamphetamine.

ee. 3,4,5-Trimethoxyamphetamine.

ff. 4-APB (4-(2-Aminopropyl)benzofuran).

gg. 5-APB (5-(2-Aminopropyl)benzofuran).

hh. 6-APB (6-(2-Aminopropyl)benzofuran).

ii. 7-APB (7-(2-Aminopropyl)benzofuran).

jj. 4-APDB (4-(2-Aminopropyl)-2,3-dihydrobenzofuran).

kk. 5-APDB (5-(2-Aminopropyl)-2,3-dihydrobenzofuran).

ll. 6-APDB (6-(2-Aminopropyl)-2,3-dihydrobenzofuran).

mm. 7-APDB (7-(2-Aminopropyl)-2,3-dihydrobenzofuran).

nn. 4-MAPB (4-(2-Methylaminopropyl)benzofuran).

oo. 5-MAPB (5-(2-Methylaminopropyl)benzofuran).

pp. 6-MAPB (6-(2-Methylaminopropyl)benzofuran).

qq. 7-MAPB (7-(2-Methylaminopropyl)benzofuran).

rr. 5-EAPB (5-(2-Ethylaminopropyl)benzofuran).

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

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).

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-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).
l. 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 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-(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)64. ~~s. 893.03(1)(a)63.~~;

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

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

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

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

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 (a) of subsection (2) of section 893.131, Florida Statutes, is amended to read:

893.131 Distribution of controlled substances resulting in overdose or serious bodily injury.—

(2)(a) Except as provided in paragraph (b), a person 18 years of age or older who unlawfully distributes:

1. Heroin, as described in s. 893.03(1)(b)11.;

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

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

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

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

6. Fentanyl derivatives, as described in s. 893.03(1)(a)64. ~~s. 893.03(1)(a)63.~~;

7. A controlled substance analog, as described in s. 893.0356, of any substance specified in subparagraphs 1.-6.; or

8. A mixture containing any substance specified in

subparagraphs 1.-7.,

and an overdose or serious bodily injury of the user results, commits a felony of the second degree, punishable as provided in s. 775.082, s. 775.083, or s. 775.084, when such substance or mixture is proven to have caused or been a substantial factor in causing the overdose or serious bodily injury of the user.

Section 4. 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 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 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.

4.a. 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:

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

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

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

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

(V) A fentanyl derivative, as described in s. 893.03(1)(a)64. ~~s. 893.03(1)(a)63.~~;

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

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

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

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

(I) Is 4 grams or more, but less than 14 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 $50,000.

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

(III) Is 28 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.

c. A person 18 years of age or older who violates sub-subparagraph a. by knowingly selling or delivering to a minor at least 4 grams of a substance or mixture listed in sub-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 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 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 $50,000.

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 5. Except as otherwise expressly provided in this act and except for this section, which shall take effect upon this act becoming a law, this act shall take effect October 1, 2026.