

1                                   A bill to be entitled  
 2           An act relating to pharmaceutical products containing  
 3           cannabis; amending s. 893.03, F.S.; excluding from  
 4           Schedule I cannabis if it is contained within a  
 5           pharmaceutical product approved by the United States  
 6           Food and Drug Administration; removing provisions  
 7           concerning the scheduling of certain drug products  
 8           containing cannabidiol; amending s. 893.02, F.S.;  
 9           conforming provisions to changes made by the act;  
 10          providing an effective date.

11

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

13

14           Section 1. Paragraph (c) of subsection (1) and paragraph  
 15           (d) of subsection (5) of section 893.03, Florida Statutes, are  
 16           amended to read:

17           893.03 Standards and schedules.—The substances enumerated  
 18           in this section are controlled by this chapter. The controlled  
 19           substances listed or to be listed in Schedules I, II, III, IV,  
 20           and V are included by whatever official, common, usual,  
 21           chemical, trade name, or class designated. The provisions of  
 22           this section shall not be construed to include within any of the  
 23           schedules contained in this section any excluded drugs listed  
 24           within the purview of 21 C.F.R. s. 1308.22, styled "Excluded  
 25           Substances"; 21 C.F.R. s. 1308.24, styled "Exempt Chemical

26 Preparations"; 21 C.F.R. s. 1308.32, styled "Exempted  
 27 Prescription Products"; or 21 C.F.R. s. 1308.34, styled "Exempt  
 28 Anabolic Steroid Products."

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

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

- 44 1. Alpha-Ethyltryptamine.
- 45 2. 4-Methylaminorex (2-Amino-4-methyl-5-phenyl-2-  
 46 oxazoline).
- 47 3. Aminorex (2-Amino-5-phenyl-2-oxazoline).
- 48 4. DOB (4-Bromo-2,5-dimethoxyamphetamine).
- 49 5. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
- 50 6. Bufotenine.

- 51           7. Cannabis, except if it is contained within a
- 52 pharmaceutical product approved by the United States Food and
- 53 Drug Administration.
- 54           8. Cathinone.
- 55           9. DET (Diethyltryptamine).
- 56           10. 2,5-Dimethoxyamphetamine.
- 57           11. DOET (4-Ethyl-2,5-Dimethoxyamphetamine).
- 58           12. DMT (Dimethyltryptamine).
- 59           13. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine
- 60 analog of phencyclidine).
- 61           14. JB-318 (N-Ethyl-3-piperidyl benzilate).
- 62           15. N-Ethylamphetamine.
- 63           16. Fenethylamine.
- 64           17. 3,4-Methylenedioxy-N-hydroxyamphetamine.
- 65           18. Ibogaine.
- 66           19. LSD (Lysergic acid diethylamide).
- 67           20. Mescaline.
- 68           21. Methcathinone.
- 69           22. 5-Methoxy-3,4-methylenedioxyamphetamine.
- 70           23. PMA (4-Methoxyamphetamine).
- 71           24. PMMA (4-Methoxymethamphetamine).
- 72           25. DOM (4-Methyl-2,5-dimethoxyamphetamine).
- 73           26. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
- 74           27. MDA (3,4-Methylenedioxyamphetamine).
- 75           28. JB-336 (N-Methyl-3-piperidyl benzilate).

- 76 |           29. N,N-Dimethylamphetamine.
- 77 |           30. Parahexyl.
- 78 |           31. Peyote.
- 79 |           32. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine
- 80 | analog of phencyclidine).
- 81 |           33. Psilocybin.
- 82 |           34. Psilocyn.
- 83 |           35. *Salvia divinorum*, except for any drug product approved
- 84 | by the United States Food and Drug Administration which contains
- 85 | *Salvia divinorum* or its isomers, esters, ethers, salts, and
- 86 | salts of isomers, esters, and ethers, if the existence of such
- 87 | isomers, esters, ethers, and salts is possible within the
- 88 | specific chemical designation.
- 89 |           36. Salvinorin A, except for any drug product approved by
- 90 | the United States Food and Drug Administration which contains
- 91 | Salvinorin A or its isomers, esters, ethers, salts, and salts of
- 92 | isomers, esters, and ethers, if the existence of such isomers,
- 93 | esters, ethers, and salts is possible within the specific
- 94 | chemical designation.
- 95 |           37. Xylazine.
- 96 |           38. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine)
- 97 | (Thiophene analog of phencyclidine).
- 98 |           39. 3,4,5-Trimethoxyamphetamine.
- 99 |           40. Methylone (3,4-Methylenedioxymethcathinone).
- 100 |           41. MDPV (3,4-Methylenedioxypyrovalerone).

- 101 42. Methylnmethcathinone.
- 102 43. Methoxymethcathinone.
- 103 44. Fluoromethcathinone.
- 104 45. Methylethcathinone.
- 105 46. CP 47,497 (2-(3-Hydroxycyclohexyl)-5-(2-methyloctan-2-
- 106 yl)phenol) and its dimethyloctyl (C8) homologue.
- 107 47. HU-210 [(6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-
- 108 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-
- 109 ol].
- 110 48. JWH-018 (1-Pentyl-3-(1-naphthoyl)indole).
- 111 49. JWH-073 (1-Butyl-3-(1-naphthoyl)indole).
- 112 50. JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-
- 113 naphthoyl)indole).
- 114 51. BZP (Benzylpiperazine).
- 115 52. Fluorophenylpiperazine.
- 116 53. Methylphenylpiperazine.
- 117 54. Chlorophenylpiperazine.
- 118 55. Methoxyphenylpiperazine.
- 119 56. DBZP (1,4-Dibenzylpiperazine).
- 120 57. TFMPP (Trifluoromethylphenylpiperazine).
- 121 58. MBDB (Methylbenzodioxolylbutanamine) or (3,4-
- 122 Methylenedioxy-N-methylbutanamine).
- 123 59. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
- 124 60. 5-Hydroxy-N-methyltryptamine.
- 125 61. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).

- 126 | 62. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
- 127 | 63. Methyltryptamine.
- 128 | 64. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).
- 129 | 65. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
- 130 | 66. Tyramine (4-Hydroxyphenethylamine).
- 131 | 67. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
- 132 | 68. DiPT (N,N-Diisopropyltryptamine).
- 133 | 69. DPT (N,N-Dipropyltryptamine).
- 134 | 70. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
- 135 | 71. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
- 136 | 72. DOI (4-Iodo-2,5-dimethoxyamphetamine).
- 137 | 73. DOC (4-Chloro-2,5-dimethoxyamphetamine).
- 138 | 74. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
- 139 | 75. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
- 140 | 76. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
- 141 | 77. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
- 142 | 78. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
- 143 | 79. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
- 144 | 80. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
- 145 | 81. Butylone (3,4-Methylenedioxy-alpha-
- 146 | methylaminobutyrophenone).
- 147 | 82. Ethcathinone.
- 148 | 83. Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
- 149 | 84. Naphyrone (Naphthylpyrovalerone).
- 150 | 85. Dimethylone (3,4-Methylenedioxy-N,N-

- 151 dimethylcathinone).
- 152 86. 3,4-Methylenedioxy-N,N-diethylcathinone.
- 153 87. 3,4-Methylenedioxy-propiofenone.
- 154 88. 3,4-Methylenedioxy-alpha-bromopropiofenone.
- 155 89. 3,4-Methylenedioxy-propiofenone-2-oxime.
- 156 90. 3,4-Methylenedioxy-N-acetylcathinone.
- 157 91. 3,4-Methylenedioxy-N-acetylmethcathinone.
- 158 92. 3,4-Methylenedioxy-N-acetylethcathinone.
- 159 93. Bromomethcathinone.
- 160 94. Buphedrone (alpha-Methylamino-butyrophenone).
- 161 95. Eutylone (3,4-Methylenedioxy-alpha-
- 162 ethylaminobutyrophenone).
- 163 96. Dimethylcathinone.
- 164 97. Dimethylmethcathinone.
- 165 98. Pentylone (3,4-Methylenedioxy-alpha-
- 166 methylaminovalerophenone).
- 167 99. MDPMP (3,4-Methylenedioxy-alpha-
- 168 pyrrolidinopropiofenone).
- 169 100. MDPBP (3,4-Methylenedioxy-alpha-
- 170 pyrrolidinobutyrophenone).
- 171 101. MOPMP (Methoxy-alpha-pyrrolidinopropiofenone).
- 172 102. MPHP (Methyl-alpha-pyrrolidinohexanophenone).
- 173 103. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP
- 174 (Benocyclidine).
- 175 104. F-MABP (Fluoromethylaminobutyrophenone).

- 176 | 105. MeO-PBP (Methoxypyrrolidinobutyrophenone).
- 177 | 106. Et-PBP (Ethylpyrrolidinobutyrophenone).
- 178 | 107. 3-Me-4-MeO-MCAT (3-Methyl-4-Methoxymethcathinone).
- 179 | 108. Me-EABP (Methylethylaminobutyrophenone).
- 180 | 109. Etizolam.
- 181 | 110. PPP (Pyrrolidinopropiophenone).
- 182 | 111. PBP (Pyrrolidinobutyrophenone).
- 183 | 112. PVP (Pyrrolidinovalerophenone) or
- 184 | (Pyrrolidinopentiophenone).
- 185 | 113. MPPP (Methyl-alpha-pyrrolidinopropiophenone).
- 186 | 114. JWH-007 (1-Pentyl-2-methyl-3-(1-naphthoyl)indole).
- 187 | 115. JWH-015 (1-Propyl-2-methyl-3-(1-naphthoyl)indole).
- 188 | 116. JWH-019 (1-Hexyl-3-(1-naphthoyl)indole).
- 189 | 117. JWH-020 (1-Heptyl-3-(1-naphthoyl)indole).
- 190 | 118. JWH-072 (1-Propyl-3-(1-naphthoyl)indole).
- 191 | 119. JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl)indole).
- 192 | 120. JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
- 193 | 121. JWH-133 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-
- 194 | methylpentan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).
- 195 | 122. JWH-175 (1-Pentyl-3-(1-naphthylmethyl)indole).
- 196 | 123. JWH-201 (1-Pentyl-3-(4-methoxyphenylacetyl)indole).
- 197 | 124. JWH-203 (1-Pentyl-3-(2-chlorophenylacetyl)indole).
- 198 | 125. JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).
- 199 | 126. JWH-250 (1-Pentyl-3-(2-methoxyphenylacetyl)indole).
- 200 | 127. JWH-251 (1-Pentyl-3-(2-methylphenylacetyl)indole).



- 201 128. JWH-302 (1-Pentyl-3-(3-methoxyphenylacetyl)indole).
- 202 129. JWH-398 (1-Pentyl-3-(4-chloro-1-naphthoyl)indole).
- 203 130. HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-
- 204 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-
- 205 ol).
- 206 131. HU-308 ([ (1R,2R,5R)-2-[2,6-Dimethoxy-4-(2-
- 207 methyloctan-2-yl)phenyl]-7,7-dimethyl-4-bicyclo[3.1.1]hept-3-
- 208 enyl] methanol).
- 209 132. HU-331 (3-Hydroxy-2-[(1R,6R)-3-methyl-6-(1-
- 210 methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-2,5-cyclohexadiene-
- 211 1,4-dione).
- 212 133. CB-13 (4-Pentyloxy-1-(1-naphthoyl)naphthalene).
- 213 134. CB-25 (N-Cyclopropyl-11-(3-hydroxy-5-pentylphenoxy)-
- 214 undecanamide).
- 215 135. CB-52 (N-Cyclopropyl-11-(2-hexyl-5-hydroxyphenoxy)-
- 216 undecanamide).
- 217 136. CP 55,940 (2-[3-Hydroxy-6-propanol-cyclohexyl]-5-(2-
- 218 methyloctan-2-yl)phenol).
- 219 137. AM-694 (1-(5-Fluoropentyl)-3-(2-iodobenzoyl)indole).
- 220 138. AM-2201 (1-(5-Fluoropentyl)-3-(1-naphthoyl)indole).
- 221 139. RCS-4 (1-Pentyl-3-(4-methoxybenzoyl)indole).
- 222 140. RCS-8 (1-(2-Cyclohexylethyl)-3-(2-
- 223 methoxyphenylacetyl)indole).
- 224 141. WIN55,212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-
- 225 morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-

226 | naphthalenylmethanone) .  
227 |       142. WIN55,212-3 ([ (3S)-2,3-Dihydro-5-methyl-3-(4-  
228 | morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-  
229 | naphthalenylmethanone) .  
230 |       143. Pentedrone (alpha-Methylaminovalerophenone) .  
231 |       144. Fluoroamphetamine .  
232 |       145. Fluoromethamphetamine .  
233 |       146. Methoxetamine .  
234 |       147. Methiopropamine .  
235 |       148. Methylbuphedrone (Methyl-alpha-  
236 | methylaminobutyrophenone) .  
237 |       149. APB ((2-Aminopropyl)benzofuran) .  
238 |       150. APDB ((2-Aminopropyl)-2,3-dihydrobenzofuran) .  
239 |       151. UR-144 (1-Pentyl-3-(2,2,3,3-  
240 | tetramethylcyclopropanoyl)indole) .  
241 |       152. XLR11 (1-(5-Fluoropentyl)-3-(2,2,3,3-  
242 | tetramethylcyclopropanoyl)indole) .  
243 |       153. Chloro UR-144 (1-(Chloropentyl)-3-(2,2,3,3-  
244 | tetramethylcyclopropanoyl)indole) .  
245 |       154. AKB48 (N-Adamant-1-yl 1-pentylindazole-3-  
246 | carboxamide) .  
247 |       155. AM-2233(1-[ (N-Methyl-2-piperidinyl)methyl]-3-(2-  
248 | iodobenzoyl)indole) .  
249 |       156. STS-135 (N-Adamant-1-yl 1-(5-fluoropentyl)indole-3-  
250 | carboxamide) .

- 251           157.   URB-597 ((3'-(Aminocarbonyl)[1,1'-biphenyl]-3-yl)-  
 252   cyclohexylcarbamate).
- 253           158.   URB-602 ([1,1'-Biphenyl]-3-yl-carbamic acid,  
 254   cyclohexyl ester).
- 255           159.   URB-754 (6-Methyl-2-[(4-methylphenyl)amino]-1-  
 256   benzoxazin-4-one).
- 257           160.   2C-D (4-Methyl-2,5-dimethoxyphenethylamine).
- 258           161.   2C-H (2,5-Dimethoxyphenethylamine).
- 259           162.   2C-N (4-Nitro-2,5-dimethoxyphenethylamine).
- 260           163.   2C-P (4-(n)-Propyl-2,5-dimethoxyphenethylamine).
- 261           164.   25I-NBOMe (4-Iodo-2,5-dimethoxy-[N-(2-  
 262   methoxybenzyl)]phenethylamine).
- 263           165.   MDMA (3,4-Methylenedioxyamphetamine).
- 264           166.   PB-22 (8-Quinoliny 1-pentylindole-3-carboxylate).
- 265           167.   Fluoro PB-22 (8-Quinoliny 1-(fluoropentyl)indole-3-  
 266   carboxylate).
- 267           168.   BB-22 (8-Quinoliny 1-(cyclohexylmethyl)indole-3-  
 268   carboxylate).
- 269           169.   Fluoro AKB48 (N-Adamant-1-yl 1-  
 270   (fluoropentyl)indazole-3-carboxamide).
- 271           170.   AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-  
 272   pentylindazole-3-carboxamide).
- 273           171.   AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-  
 274   (4-fluorobenzyl)indazole-3-carboxamide).
- 275           172.   ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-

276 1-pentylindazole-3-carboxamide) .  
277 173. Fluoro ADBICA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-  
278 yl)-1-(fluoropentyl)indole-3-carboxamide) .  
279 174. 25B-NBOMe (4-Bromo-2,5-dimethoxy-[N-(2-  
280 methoxybenzyl)]phenethylamine) .  
281 175. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-  
282 methoxybenzyl)]phenethylamine) .  
283 176. AB-CHMINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-  
284 (cyclohexylmethyl)indazole-3-carboxamide) .  
285 177. FUB-PB-22 (8-Quinoliny 1-(4-fluorobenzyl)indole-3-  
286 carboxylate) .  
287 178. Fluoro-NNEI (N-Naphthalen-1-yl 1-  
288 (fluoropentyl)indole-3-carboxamide) .  
289 179. Fluoro-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-  
290 (fluoropentyl)indazole-3-carboxamide) .  
291 180. THJ-2201 (1-(5-Fluoropentyl)-3-(1-  
292 naphthoyl)indazole) .  
293 181. AM-855 ((4aR,12bR)-8-Hexyl-2,5,5-trimethyl-  
294 1,4,4a,8,9,10,11,12b-octahydronaphtho[3,2-c]isochromen-12-ol) .  
295 182. AM-905 ((6aR,9R,10aR)-3-[(E)-Hept-1-enyl]-9-  
296 (hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-  
297 hexahydrobenzo[c]chromen-1-ol) .  
298 183. AM-906 ((6aR,9R,10aR)-3-[(Z)-Hept-1-enyl]-9-  
299 (hydroxymethyl)-6,6-dimethyl-6a,7,8,9,10,10a-  
300 hexahydrobenzo[c]chromen-1-ol) .

301 184. AM-2389 ((6aR, 9R, 10aR)-3-(1-Hexyl-cyclobut-1-yl)-  
 302 6a, 7, 8, 9, 10, 10a-hexahydro-6, 6-dimethyl-6H-dibenzo[b, d]pyran-1, 9  
 303 diol).

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

307 186. HU-336 ((6aR, 10aR)-6, 6, 9-Trimethyl-3-pentyl-  
 308 6a, 7, 10, 10a-tetrahydro-1H-benzo[c]chromene-1, 4(6H)-dione).

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

310 188. 5-IT (2-(1H-Indol-5-yl)-1-methyl-ethylamine).

311 189. 6-IT (2-(1H-Indol-6-yl)-1-methyl-ethylamine).

312 190. Synthetic Cannabinoids.—Unless specifically excepted  
 313 or unless listed in another schedule or contained within a  
 314 pharmaceutical product approved by the United States Food and  
 315 Drug Administration, any material, compound, mixture, or  
 316 preparation that contains any quantity of a synthetic  
 317 cannabinoid found to be in any of the following chemical class  
 318 descriptions, or homologues, nitrogen-heterocyclic analogs,  
 319 isomers (including optical, positional, or geometric), esters,  
 320 ethers, salts, and salts of homologues, nitrogen-heterocyclic  
 321 analogs, isomers, esters, or ethers, whenever the existence of  
 322 such homologues, nitrogen-heterocyclic analogs, isomers, esters,  
 323 ethers, salts, and salts of isomers, esters, or ethers is  
 324 possible within the specific chemical class or designation.  
 325 Since nomenclature of these synthetically produced cannabinoids

326 is not internationally standardized and may continually evolve,  
 327 these structures or the compounds of these structures shall be  
 328 included under this subparagraph, regardless of their specific  
 329 numerical designation of atomic positions covered, if it can be  
 330 determined through a recognized method of scientific testing or  
 331 analysis that the substance contains properties that fit within  
 332 one or more of the following categories:

333 a. Tetrahydrocannabinols.—Any tetrahydrocannabinols  
 334 naturally contained in a plant of the genus *Cannabis*, the  
 335 synthetic equivalents of the substances contained in the plant  
 336 or in the resinous extracts of the genus *Cannabis*, or synthetic  
 337 substances, derivatives, and their isomers with similar chemical  
 338 structure and pharmacological activity, including, but not  
 339 limited to, Delta 9 tetrahydrocannabinols and their optical  
 340 isomers, Delta 8 tetrahydrocannabinols and their optical  
 341 isomers, Delta 6a,10a tetrahydrocannabinols and their optical  
 342 isomers, or any compound containing a tetrahydrobenzo[c]chromene  
 343 structure with substitution at either or both the 3-position or  
 344 9-position, with or without substitution at the 1-position with  
 345 hydroxyl or alkoxy groups, including, but not limited to:

- 346 (I) Tetrahydrocannabinol.
- 347 (II) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-  
 348 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-  
 349 ol).
- 350 (III) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-

351 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-  
352 ol).

353 (IV) JWH-051 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-  
354 (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

355 (V) JWH-133 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-methylpentan-  
356 2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

357 (VI) JWH-057 ((6aR,10aR)-6,6,9-Trimethyl-3-(2-methyloctan-  
358 2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromene).

359 (VII) JWH-359 ((6aR,10aR)-1-Methoxy-6,6,9-trimethyl-3-  
360 (2,3-dimethylpentan-2-yl)-6a,7,10,10a-  
361 tetrahydrobenzo[c]chromene).

362 (VIII) AM-087 ((6aR,10aR)-3-(2-Methyl-6-bromohex-2-yl)-  
363 6,6,9-trimethyl-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol).

364 (IX) AM-411 ((6aR,10aR)-3-(1-Adamantyl)-6,6,9-trimethyl-  
365 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol).

366 (X) Parahexyl.

367 b. Naphthoylindoles, Naphthoylindazoles,  
368 Naphthoylcarbazoles, Naphthylmethylindoles,  
369 Naphthylmethylindazoles, and Naphthylmethylcarbazoles.—Any  
370 compound containing a naphthoylindole, naphthoylindazole,  
371 naphthoylcarbazole, naphthylmethylindole,  
372 naphthylmethylindazole, or naphthylmethylcarbazole structure,  
373 with or without substitution on the indole, indazole, or  
374 carbazole ring to any extent, whether or not substituted on the  
375 naphthyl ring to any extent, including, but not limited to:

- 376 (I) JWH-007 (1-Pentyl-2-methyl-3-(1-naphthoyl) indole) .
- 377 (II) JWH-011 (1-(1-Methylhexyl)-2-methyl-3-(1-
- 378 naphthoyl) indole) .
- 379 (III) JWH-015 (1-Propyl-2-methyl-3-(1-naphthoyl) indole) .
- 380 (IV) JWH-016 (1-Butyl-2-methyl-3-(1-naphthoyl) indole) .
- 381 (V) JWH-018 (1-Pentyl-3-(1-naphthoyl) indole) .
- 382 (VI) JWH-019 (1-Hexyl-3-(1-naphthoyl) indole) .
- 383 (VII) JWH-020 (1-Heptyl-3-(1-naphthoyl) indole) .
- 384 (VIII) JWH-022 (1-(4-Pentenyl)-3-(1-naphthoyl) indole) .
- 385 (IX) JWH-071 (1-Ethyl-3-(1-naphthoyl) indole) .
- 386 (X) JWH-072 (1-Propyl-3-(1-naphthoyl) indole) .
- 387 (XI) JWH-073 (1-Butyl-3-(1-naphthoyl) indole) .
- 388 (XII) JWH-080 (1-Butyl-3-(4-methoxy-1-naphthoyl) indole) .
- 389 (XIII) JWH-081 (1-Pentyl-3-(4-methoxy-1-naphthoyl) indole) .
- 390 (XIV) JWH-098 (1-Pentyl-2-methyl-3-(4-methoxy-1-
- 391 naphthoyl) indole) .
- 392 (XV) JWH-116 (1-Pentyl-2-ethyl-3-(1-naphthoyl) indole) .
- 393 (XVI) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl) indole) .
- 394 (XVII) JWH-149 (1-Pentyl-2-methyl-3-(4-methyl-1-
- 395 naphthoyl) indole) .
- 396 (XVIII) JWH-164 (1-Pentyl-3-(7-methoxy-1-
- 397 naphthoyl) indole) .
- 398 (XIX) JWH-175 (1-Pentyl-3-(1-naphthylmethyl) indole) .
- 399 (XX) JWH-180 (1-Propyl-3-(4-propyl-1-naphthoyl) indole) .
- 400 (XXI) JWH-182 (1-Pentyl-3-(4-propyl-1-naphthoyl) indole) .



401 (XXII) JWH-184 (1-Pentyl-3-[(4-methyl)-1-  
 402 naphthylmethyl]indole).  
 403 (XXIII) JWH-193 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methyl-1-  
 404 naphthoyl)indole).  
 405 (XXIV) JWH-198 (1-[2-(4-Morpholinyl)ethyl]-3-(4-methoxy-1-  
 406 naphthoyl)indole).  
 407 (XXV) JWH-200 (1-[2-(4-Morpholinyl)ethyl]-3-(1-  
 408 naphthoyl)indole).  
 409 (XXVI) JWH-210 (1-Pentyl-3-(4-ethyl-1-naphthoyl)indole).  
 410 (XXVII) JWH-387 (1-Pentyl-3-(4-bromo-1-naphthoyl)indole).  
 411 (XXVIII) JWH-398 (1-Pentyl-3-(4-chloro-1-  
 412 naphthoyl)indole).  
 413 (XXIX) JWH-412 (1-Pentyl-3-(4-fluoro-1-naphthoyl)indole).  
 414 (XXX) JWH-424 (1-Pentyl-3-(8-bromo-1-naphthoyl)indole).  
 415 (XXXI) AM-1220 (1-[(1-Methyl-2-piperidinyl)methyl]-3-(1-  
 416 naphthoyl)indole).  
 417 (XXXII) AM-1235 (1-(5-Fluoropentyl)-6-nitro-3-(1-  
 418 naphthoyl)indole).  
 419 (XXXIII) AM-2201 (1-(5-Fluoropentyl)-3-(1-  
 420 naphthoyl)indole).  
 421 (XXXIV) Chloro JWH-018 (1-(Chloropentyl)-3-(1-  
 422 naphthoyl)indole).  
 423 (XXXV) Bromo JWH-018 (1-(Bromopentyl)-3-(1-  
 424 naphthoyl)indole).  
 425 (XXXVI) AM-2232 (1-(4-Cyanobutyl)-3-(1-naphthoyl)indole).

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

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

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

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

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

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

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

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

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

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

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

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

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

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

450 (IX) JWH-369 (1-Pentyl-5-(2-chlorophenyl)-3-(1-

451 naphthoyl)pyrrole).

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

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

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

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

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

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

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

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

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

471 (VII) Cannabipiperidiethanone.

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

474 f. Cyclohexylphenols.—Any compound containing a  
475 cyclohexylphenol structure, with or without substitution at the

476 5-position of the phenolic ring to any extent, whether or not  
477 substituted on the cyclohexyl ring to any extent, including, but  
478 not limited to:

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

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

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

485 g. Benzoylindoles and Benzoylindazoles.—Any compound  
486 containing a benzoylindole or benzoylindazole structure, with or  
487 without substitution on the indole or indazole ring to any  
488 extent, whether or not substituted on the phenyl ring to any  
489 extent, including, but not limited to:

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

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

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

494 (IV) Pravadoline (1-[2-(4-Morpholinyl)ethyl]-2-methyl-3-  
495 (4-methoxybenzoyl)indole).

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

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

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

501 (VIII) AM-630 (1-[2-(4-Morpholinyl)ethyl]-2-methyl-6-iodo-  
 502 3-(4-methoxybenzoyl)indole).

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

548 j. Quinolinyndolecarboxylates,  
 549 Quinolinyndazolecarboxylates, Quinolinyndolecarboxamides,  
 550 and Quinolinyndazolecarboxamides.—Any compound containing a

551 quinolinylindole carboxylate, quinolinylindazole carboxylate,  
552 isoquinolinylindole carboxylate, isoquinolinylindazole  
553 carboxylate, quinolinylindole carboxamide, quinolinylindazole  
554 carboxamide, isoquinolinylindole carboxamide, or  
555 isoquinolinylindazole carboxamide structure, with or without  
556 substitution on the indole or indazole ring to any extent,  
557 whether or not substituted on the quinoline or isoquinoline ring  
558 to any extent, including, but not limited to:

559 (I) PB-22 (8-Quinolinyl 1-pentylindole-3-carboxylate).

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

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

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

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

567 (VI) Fluoro NPB-22 (8-Quinolinyl 1-(fluoropentyl)indazole-  
568 3-carboxylate).

569 (VII) FUB-NPB-22 (8-Quinolinyl 1-(4-fluorobenzyl)indazole-  
570 3-carboxylate).

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

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

574 k. Naphthylindolecarboxylates and  
575 Naphthylindazolecarboxylates.—Any compound containing a

576 naphthylindole carboxylate or naphthylindazole carboxylate  
 577 structure, with or without substitution on the indole or  
 578 indazole ring to any extent, whether or not substituted on the  
 579 naphthyl ring to any extent, including, but not limited to:

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

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

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

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

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

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

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

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

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



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

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

605 m. Alkylcarbonyl indole carboxamides, Alkylcarbonyl  
 606 indazole carboxamides, Alkylcarbonyl indole carboxylates, and  
 607 Alkylcarbonyl indazole carboxylates.—Any compound containing an  
 608 alkylcarbonyl group, including 1-amino-3-methyl-1-oxobutan-2-yl,  
 609 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-amino-1-oxo-3-  
 610 phenylpropan-2-yl, 1-methoxy-1-oxo-3-phenylpropan-2-yl, with an  
 611 indole carboxamide, indazole carboxamide, indole carboxylate, or  
 612 indazole carboxylate, with or without substitution on the indole  
 613 or indazole ring to any extent, whether or not substituted on  
 614 the alkylcarbonyl group to any extent, including, but not  
 615 limited to:

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

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

620 (III) Fluoro ABICA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-  
 621 1-(fluoropentyl)indole-3-carboxamide).

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

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

- 626 (VI) ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-  
627 1-pentylindazole-3-carboxamide).
- 628 (VII) Fluoro ADB-PINACA (N-(1-Amino-3,3-dimethyl-1-  
629 oxobutan-2-yl)-1-(fluoropentyl)indazole-3-carboxamide).
- 630 (VIII) AB-FUBINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-  
631 1-(4-fluorobenzyl)indazole-3-carboxamide).
- 632 (IX) ADB-FUBINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-  
633 yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).
- 634 (X) AB-CHMINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-  
635 (cyclohexylmethyl)indazole-3-carboxamide).
- 636 (XI) MA-CHMINACA (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-  
637 1-(cyclohexylmethyl)indazole-3-carboxamide).
- 638 (XII) MAB-CHMINACA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-  
639 yl)-1-(cyclohexylmethyl)indazole-3-carboxamide).
- 640 (XIII) AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-  
641 pentylindazole-3-carboxamide).
- 642 (XIV) Fluoro-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-  
643 1-(fluoropentyl)indazole-3-carboxamide).
- 644 (XV) FUB-AMB (N-(1-Methoxy-3-methyl-1-oxobutan-2-yl)-1-(4-  
645 fluorobenzyl)indazole-3-carboxamide).
- 646 (XVI) MDMB-CHMINACA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-  
647 2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide).
- 648 (XVII) MDMB-FUBINACA (N-(1-Methoxy-3,3-dimethyl-1-  
649 oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide).
- 650 (XVIII) MDMB-CHMICA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-

651 2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide).

652 (XIX) PX-1 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(5-

653 fluoropentyl)indole-3-carboxamide).

654 (XX) PX-2 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(5-

655 fluoropentyl)indazole-3-carboxamide).

656 (XXI) PX-3 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-

657 (cyclohexylmethyl)indazole-3-carboxamide).

658 (XXII) PX-4 (N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(4-

659 fluorobenzyl)indazole-3-carboxamide).

660 (XXIII) MO-CHMINACA (N-(1-Methoxy-3,3-dimethyl-1-oxobutan-

661 2-yl)-1-(cyclohexylmethyl)indazole-3-carboxylate).

662 n. Cumylindolecarboxamides and Cumylindazolecarboxamides.-

663 Any compound containing a N-(2-phenylpropan-2-yl) indole

664 carboxamide or N-(2-phenylpropan-2-yl) indazole carboxamide

665 structure, with or without substitution on the indole or

666 indazole ring to any extent, whether or not substituted on the

667 phenyl ring of the cumyl group to any extent, including, but not

668 limited to:

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

670 carboxamide).

671 (II) Fluoro CUMYL-PICA (N-(2-Phenylpropan-2-yl)-1-

672 (fluoropentyl)indole-3-carboxamide).

673 o. Other Synthetic Cannabinoids.-Any material, compound,

674 mixture, or preparation that contains any quantity of a

675 Synthetic Cannabinoid, as described in sub-subparagraphs a.-n.:

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

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

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

688 191. Substituted Cathinones.—Unless specifically excepted,  
 689 listed in another schedule, or contained within a pharmaceutical  
 690 product approved by the United States Food and Drug  
 691 Administration, any material, compound, mixture, or preparation,  
 692 including its salts, isomers, esters, or ethers, and salts of  
 693 isomers, esters, or ethers, whenever the existence of such salts  
 694 is possible within any of the following specific chemical  
 695 designations:

696 a. Any compound containing a 2-amino-1-phenyl-1-propanone  
 697 structure;

698 b. Any compound containing a 2-amino-1-naphthyl-1-  
 699 propanone structure; or

700 c. Any compound containing a 2-amino-1-thiophenyl-1-

701 propanone structure,

702

703 whether or not the compound is further modified:

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

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

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

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

- 717 (A) Methcathinone.
- 718 (B) Ethcathinone.
- 719 (C) Methylone (3,4-Methylenedioxy-methcathinone).
- 720 (D) 2,3-Methylenedioxy-methcathinone.
- 721 (E) MDPV (3,4-Methylenedioxy-pyrovalerone).
- 722 (F) Methylmethcathinone.
- 723 (G) Methoxymethcathinone.
- 724 (H) Fluoromethcathinone.
- 725 (I) Methylethcathinone.

- 726 (J) Butylone (3,4-Methylenedioxy-alpha-
- 727 methylaminobutyrophenone).
- 728 (K) Ethylone (3,4-Methylenedioxy-N-ethylcathinone).
- 729 (L) BMDP (3,4-Methylenedioxy-N-benzylcathinone).
- 730 (M) Naphyrone (Naphthylpyrovalerone).
- 731 (N) Bromomethcathinone.
- 732 (O) Buphedrone (alpha-Methylaminobutyrophenone).
- 733 (P) Eutylone (3,4-Methylenedioxy-alpha-
- 734 ethylaminobutyrophenone).
- 735 (Q) Dimethylcathinone.
- 736 (R) Dimethylmethcathinone.
- 737 (S) Pentylone (3,4-Methylenedioxy-alpha-
- 738 methylaminovalerophenone).
- 739 (T) Pentedrone (alpha-Methylaminovalerophenone).
- 740 (U) MDPPP (3,4-Methylenedioxy-alpha-
- 741 pyrrolidinopropiophenone).
- 742 (V) MDPBP (3,4-Methylenedioxy-alpha-
- 743 pyrrolidinobutyrophenone).
- 744 (W) MPPP (Methyl-alpha-pyrrolidinopropiophenone).
- 745 (X) PPP (Pyrrolidinopropiophenone).
- 746 (Y) PVP (Pyrrolidinovalerophenone) or
- 747 (Pyrrolidinopentiophenone).
- 748 (Z) MOPPP (Methoxy-alpha-pyrrolidinopropiophenone).
- 749 (AA) MPHP (Methyl-alpha-pyrrolidinohexanophenone).
- 750 (BB) F-MABP (Fluoromethylaminobutyrophenone).

- 751 (CC) Me-EABP (Methylethylaminobutyrophenone).
- 752 (DD) PBP (Pyrrolidinobutyrophenone).
- 753 (EE) MeO-PBP (Methoxypyrrolidinobutyrophenone).
- 754 (FF) Et-PBP (Ethylpyrrolidinobutyrophenone).
- 755 (GG) 3-Me-4-MeO-MCAT (3-Methyl-4-Methoxymethcathinone).
- 756 (HH) Dimethylone (3,4-Methylenedioxy-N,N-
- 757 dimethylcathinone).
- 758 (II) 3,4-Methylenedioxy-N,N-diethylcathinone.
- 759 (JJ) 3,4-Methylenedioxy-N-acetylcathinone.
- 760 (KK) 3,4-Methylenedioxy-N-acetylmethcathinone.
- 761 (LL) 3,4-Methylenedioxy-N-acetylethcathinone.
- 762 (MM) Methylbuphedrone (Methyl-alpha-
- 763 methylaminobutyrophenone).
- 764 (NN) Methyl-alpha-methylaminohexanophenone.
- 765 (OO) N-Ethyl-N-methylcathinone.
- 766 (PP) PHP (Pyrrolidinohexanophenone).
- 767 (QQ) PV8 (Pyrrolidinoheptanophenone).
- 768 (RR) Chloromethcathinone.
- 769 (SS) 4-Bromo-2,5-dimethoxy-alpha-aminoacetophenone.
- 770 192. Substituted Phenethylamines.—Unless specifically
- 771 excepted or unless listed in another schedule, or contained
- 772 within a pharmaceutical product approved by the United States
- 773 Food and Drug Administration, any material, compound, mixture,
- 774 or preparation, including its salts, isomers, esters, or ethers,
- 775 and salts of isomers, esters, or ethers, whenever the existence

776 of such salts is possible within any of the following specific  
777 chemical designations, any compound containing a phenethylamine  
778 structure, without a beta-keto group, and without a benzyl group  
779 attached to the amine group, whether or not the compound is  
780 further modified with or without substitution on the phenyl ring  
781 to any extent with alkyl, alkylthio, nitro, alkoxy, thio,  
782 halide, fused alkylenedioxy, fused furan, fused benzofuran,  
783 fused dihydrofuran, or fused tetrahydropyran substituents,  
784 whether or not further substituted on a ring to any extent, with  
785 or without substitution at the alpha or beta position by any  
786 alkyl substituent, with or without substitution at the nitrogen  
787 atom, and with or without inclusion of the 2-amino nitrogen atom  
788 in a cyclic structure, including, but not limited to:

- 789 a. 2C-B (4-Bromo-2,5-dimethoxyphenethylamine).
- 790 b. 2C-E (4-Ethyl-2,5-dimethoxyphenethylamine).
- 791 c. 2C-T-4 (4-Isopropylthio-2,5-dimethoxyphenethylamine).
- 792 d. 2C-C (4-Chloro-2,5-dimethoxyphenethylamine).
- 793 e. 2C-T (4-Methylthio-2,5-dimethoxyphenethylamine).
- 794 f. 2C-T-2 (4-Ethylthio-2,5-dimethoxyphenethylamine).
- 795 g. 2C-T-7 (4-(n)-Propylthio-2,5-dimethoxyphenethylamine).
- 796 h. 2C-I (4-Iodo-2,5-dimethoxyphenethylamine).
- 797 i. 2C-D (4-Methyl-2,5-dimethoxyphenethylamine).
- 798 j. 2C-H (2,5-Dimethoxyphenethylamine).
- 799 k. 2C-N (4-Nitro-2,5-dimethoxyphenethylamine).
- 800 l. 2C-P (4-(n)-Propyl-2,5-dimethoxyphenethylamine).



- 801 m. MDMA (3,4-Methylenedioxyamphetamine).
- 802 n. MBDB (Methylbenzodioxolylbutanamine) or (3,4-
- 803 Methylenedioxy-N-methylbutanamine).
- 804 o. MDA (3,4-Methylenedioxyamphetamine).
- 805 p. 2,5-Dimethoxyamphetamine.
- 806 q. Fluoroamphetamine.
- 807 r. Fluoromethamphetamine.
- 808 s. MDEA (3,4-Methylenedioxy-N-ethylamphetamine).
- 809 t. DOB (4-Bromo-2,5-dimethoxyamphetamine).
- 810 u. DOC (4-Chloro-2,5-dimethoxyamphetamine).
- 811 v. DOET (4-Ethyl-2,5-dimethoxyamphetamine).
- 812 w. DOI (4-Iodo-2,5-dimethoxyamphetamine).
- 813 x. DOM (4-Methyl-2,5-dimethoxyamphetamine).
- 814 y. PMA (4-Methoxyamphetamine).
- 815 z. N-Ethylamphetamine.
- 816 aa. 3,4-Methylenedioxy-N-hydroxyamphetamine.
- 817 bb. 5-Methoxy-3,4-methylenedioxyamphetamine.
- 818 cc. PMMA (4-Methoxymethamphetamine).
- 819 dd. N,N-Dimethylamphetamine.
- 820 ee. 3,4,5-Trimethoxyamphetamine.
- 821 ff. 4-APB (4-(2-Aminopropyl)benzofuran).
- 822 gg. 5-APB (5-(2-Aminopropyl)benzofuran).
- 823 hh. 6-APB (6-(2-Aminopropyl)benzofuran).
- 824 ii. 7-APB (7-(2-Aminopropyl)benzofuran).
- 825 jj. 4-APDB (4-(2-Aminopropyl)-2,3-dihydrobenzofuran).

- 826 | kk. 5-APDB (5-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 827 | ll. 6-APDB (6-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 828 | mm. 7-APDB (7-(2-Aminopropyl)-2,3-dihydrobenzofuran).
- 829 | nn. 4-MAPB (4-(2-Methylaminopropyl)benzofuran).
- 830 | oo. 5-MAPB (5-(2-Methylaminopropyl)benzofuran).
- 831 | pp. 6-MAPB (6-(2-Methylaminopropyl)benzofuran).
- 832 | qq. 7-MAPB (7-(2-Methylaminopropyl)benzofuran).
- 833 | rr. 5-EAPB (5-(2-Ethylaminopropyl)benzofuran).
- 834 | ss. 5-MAPDB (5-(2-Methylaminopropyl)-2,3-

835 | dihydrobenzofuran),

836 |

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

842 | 193. N-Benzyl Phenethylamine Compounds.—Unless  
 843 | specifically excepted or unless listed in another schedule, or  
 844 | contained within a pharmaceutical product approved by the United  
 845 | States Food and Drug Administration, any material, compound,  
 846 | mixture, or preparation, including its salts, isomers, esters,  
 847 | or ethers, and salts of isomers, esters, or ethers, whenever the  
 848 | existence of such salts is possible within any of the following  
 849 | specific chemical designations, any compound containing a  
 850 | phenethylamine structure without a beta-keto group, with

851 substitution on the nitrogen atom of the amino group with a  
852 benzyl substituent, with or without substitution on the phenyl  
853 or benzyl ring to any extent with alkyl, alkoxy, thio,  
854 alkylthio, halide, fused alkylendioxy, fused furan, fused  
855 benzofuran, or fused tetrahydropyran substituents, whether or  
856 not further substituted on a ring to any extent, with or without  
857 substitution at the alpha position by any alkyl substituent,  
858 including, but not limited to:

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

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

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

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

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

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

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

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

875 i. 25T2-NBOMe (4-Methylthio-2,5-dimethoxy-[N-(2-

876 methoxybenzyl) ]phenethylamine) .  
 877       j. 25T4-NBOMe (4-Isopropylthio-2,5-dimethoxy-[N-(2-  
 878 methoxybenzyl) ]phenethylamine) .  
 879       k. 25T7-NBOMe (4-(n)-Propylthio-2,5-dimethoxy-[N-(2-  
 880 methoxybenzyl) ]phenethylamine) .  
 881       l. 25C-NBOMe (4-Chloro-2,5-dimethoxy-[N-(2-  
 882 methoxybenzyl) ]phenethylamine) .  
 883       m. 25C-NBOH (4-Chloro-2,5-dimethoxy-[N-(2-  
 884 hydroxybenzyl) ]phenethylamine) .  
 885       n. 25C-NBF (4-Chloro-2,5-dimethoxy-[N-(2-  
 886 fluorobenzyl) ]phenethylamine) .  
 887       o. 25C-NBMD (4-Chloro-2,5-dimethoxy-[N-(2,3-  
 888 methylenedioxybenzyl) ]phenethylamine) .  
 889       p. 25H-NBOMe (2,5-Dimethoxy-[N-(2-  
 890 methoxybenzyl) ]phenethylamine) .  
 891       q. 25H-NBOH (2,5-Dimethoxy-[N-(2-  
 892 hydroxybenzyl) ]phenethylamine) .  
 893       r. 25H-NBF (2,5-Dimethoxy-[N-(2-  
 894 fluorobenzyl) ]phenethylamine) .  
 895       s. 25D-NBOMe (4-Methyl-2,5-dimethoxy-[N-(2-  
 896 methoxybenzyl) ]phenethylamine) ,  
 897  
 898 which does not include substituted cathinones as described in  
 899 subparagraph 191.  
 900       194. Substituted Tryptamines.—Unless specifically excepted

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

- 912 a. Alpha-Ethyltryptamine.
- 913 b. Bufotenine.
- 914 c. DET (Diethyltryptamine).
- 915 d. DMT (Dimethyltryptamine).
- 916 e. MET (N-Methyl-N-ethyltryptamine).
- 917 f. DALT (N,N-Diallyltryptamine).
- 918 g. EiPT (N-Ethyl-N-isopropyltryptamine).
- 919 h. MiPT (N-Methyl-N-isopropyltryptamine).
- 920 i. 5-Hydroxy-AMT (5-Hydroxy-alpha-methyltryptamine).
- 921 j. 5-Hydroxy-N-methyltryptamine.
- 922 k. 5-MeO-MiPT (5-Methoxy-N-methyl-N-isopropyltryptamine).
- 923 l. 5-MeO-AMT (5-Methoxy-alpha-methyltryptamine).
- 924 m. Methyltryptamine.
- 925 n. 5-MeO-DMT (5-Methoxy-N,N-dimethyltryptamine).

- 926 o. 5-Me-DMT (5-Methyl-N,N-dimethyltryptamine).
- 927 p. 5-MeO-DiPT (5-Methoxy-N,N-Diisopropyltryptamine).
- 928 q. DiPT (N,N-Diisopropyltryptamine).
- 929 r. DPT (N,N-Dipropyltryptamine).
- 930 s. 4-Hydroxy-DiPT (4-Hydroxy-N,N-diisopropyltryptamine).
- 931 t. 5-MeO-DALT (5-Methoxy-N,N-Diallyltryptamine).
- 932 u. 4-AcO-DMT (4-Acetoxy-N,N-dimethyltryptamine).
- 933 v. 4-AcO-DiPT (4-Acetoxy-N,N-diisopropyltryptamine).
- 934 w. 4-Hydroxy-DET (4-Hydroxy-N,N-diethyltryptamine).
- 935 x. 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-ethyltryptamine).
- 936 y. 4-Hydroxy-MiPT (4-Hydroxy-N-methyl-N-
- 937 isopropyltryptamine).
- 938 z. Methyl-alpha-ethyltryptamine.
- 939 aa. Bromo-DALT (Bromo-N,N-diallyltryptamine),

940  
 941 which does not include tryptamine, psilocyn as described in  
 942 subparagraph 34., or psilocybin as described in subparagraph 33.

943 195. Substituted Phenylcyclohexylamines.—Unless  
 944 specifically excepted or unless listed in another schedule, or  
 945 contained within a pharmaceutical product approved by the United  
 946 States Food and Drug Administration, any material, compound,  
 947 mixture, or preparation containing a phenylcyclohexylamine  
 948 structure, with or without any substitution on the phenyl ring,  
 949 any substitution on the cyclohexyl ring, any replacement of the  
 950 phenyl ring with a thiophenyl or benzothiophenyl ring, with or

- 951 without substitution on the amine with alkyl, dialkyl, or alkoxy  
 952 substituents, inclusion of the nitrogen in a cyclic structure,  
 953 or any combination of the above, including, but not limited to:
- 954 a. BTCP (Benzothiophenylcyclohexylpiperidine) or BCP  
 955 (Benocyclidine).
  - 956 b. PCE (N-Ethyl-1-phenylcyclohexylamine) (Ethylamine analog  
 957 of phencyclidine).
  - 958 c. PCPY (N-(1-Phenylcyclohexyl)-pyrrolidine) (Pyrrolidine  
 959 analog of phencyclidine).
  - 960 d. PCPr (Phenylcyclohexylpropylamine).
  - 961 e. TCP (1-[1-(2-Thienyl)-cyclohexyl]-piperidine) (Thiophene  
 962 analog of phencyclidine).
  - 963 f. PCEEA (Phenylcyclohexyl(ethoxyethylamine)).
  - 964 g. PCMPA (Phenylcyclohexyl(methoxypropylamine)).
  - 965 h. Methoxetamine.
  - 966 i. 3-Methoxy-PCE ((3-Methoxyphenyl)cyclohexylethylamine).
  - 967 j. Bromo-PCP ((Bromophenyl)cyclohexylpiperidine).
  - 968 k. Chloro-PCP ((Chlorophenyl)cyclohexylpiperidine).
  - 969 l. Fluoro-PCP ((Fluorophenyl)cyclohexylpiperidine).
  - 970 m. Hydroxy-PCP ((Hydroxyphenyl)cyclohexylpiperidine).
  - 971 n. Methoxy-PCP ((Methoxyphenyl)cyclohexylpiperidine).
  - 972 o. Methyl-PCP ((Methylphenyl)cyclohexylpiperidine).
  - 973 p. Nitro-PCP ((Nitrophenyl)cyclohexylpiperidine).
  - 974 q. Oxo-PCP ((Oxophenyl)cyclohexylpiperidine).
  - 975 r. Amino-PCP ((Aminophenyl)cyclohexylpiperidine).

- 976 196. W-15, 4-chloro-N-[1-(2-phenylethyl)-2-  
 977 piperidinylidene]-benzenesulfonamide.
- 978 197. W-18, 4-chloro-N-[1-[2-(4-nitrophenyl)ethyl]-2-  
 979 piperidinylidene]-benzenesulfonamide.
- 980 198. AH-7921, 3,4-dichloro-N-[[1-  
 981 (dimethylamino)cyclohexyl]methyl]-benzamide.
- 982 199. U47700, trans-3,4-dichloro-N-[2-  
 983 (dimethylamino)cyclohexyl]-N-methyl-benzamide.
- 984 200. MT-45, 1-cyclohexyl-4-(1,2-diphenylethyl)-piperazine,  
 985 dihydrochloride.

986 (5) SCHEDULE V.—A substance, compound, mixture, or  
 987 preparation of a substance in Schedule V has a low potential for  
 988 abuse relative to the substances in Schedule IV and has a  
 989 currently accepted medical use in treatment in the United  
 990 States, and abuse of such compound, mixture, or preparation may  
 991 lead to limited physical or psychological dependence relative to  
 992 the substances in Schedule IV.

993 ~~(d) A drug product in finished dosage formulation that has~~  
 994 ~~been approved by the United States Food and Drug Administration~~  
 995 ~~that contains cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-~~  
 996 ~~2-cyclohexen-1-yl]-5-pentyl-1,3-benzenediol) derived from~~  
 997 ~~cannabis and no more than 0.1 percent (w/w) residual~~  
 998 ~~tetrahydrocannabinols.~~

999 Section 2. Subsection (3) of section 893.02, Florida  
 1000 Statutes, is amended to read:



1001           893.02 Definitions.—The following words and phrases as  
1002 used in this chapter shall have the following meanings, unless  
1003 the context otherwise requires:

1004           (3) "Cannabis" means all parts of any plant of the genus  
1005 Cannabis, whether growing or not; the seeds thereof; the resin  
1006 extracted from any part of the plant; and every compound,  
1007 manufacture, salt, derivative, mixture, or preparation of the  
1008 plant or its seeds or resin. The term does not include  
1009 "marijuana," as defined in s. 381.986, if manufactured,  
1010 possessed, sold, purchased, delivered, distributed, or  
1011 dispensed, in conformance with s. 381.986. The term does not  
1012 include hemp as defined in s. 581.217 or industrial hemp as  
1013 defined in s. 1004.4473. ~~The term does not include a drug~~  
1014 ~~product described in s. 893.03(5)(d).~~

1015           Section 3. This act shall take effect upon becoming a law.