

SB132 INTRODUCED



1 SB132
2 GF7H2C6-1
3 By Senator Melson
4 RFD: Healthcare
5 First Read: 06-Feb-25



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SYNOPSIS:

Existing law lists tetrahydrocannabinol (THC) as a Schedule I controlled substance, but explicitly exempts THC found in hemp.

Delta-8 THC, Delta-9 THC, and Delta-10 THC are psychoactive cannabinoids that can be made from hemp-derived cannabidiol (CBD). Under existing law, these psychoactive cannabinoids sold in Alabama may not be sold to minors and must be located in an area where minors cannot access the products.

This bill would provide that only non-psychoactive cannabinoids derived from or found in hemp are exempt from the Schedule I controlled substances list, thus classifying psychoactive cannabinoids as controlled substances.

This bill would also repeal the provision that prohibits the sale of psychoactive cannabinoids to minors.

A BILL
TO BE ENTITLED
AN ACT

Relating to controlled substances; to amend Section



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29 20-2-23, Code of Alabama 1975, to provide further for
30 tetrahydrocannabinol listed as a Schedule I controlled
31 substance; and to repeal Section 13A-12-214.4, Code of Alabama
32 1975, prohibiting the sale of psychoactive cannabinoids to
33 minors.

34 BE IT ENACTED BY THE LEGISLATURE OF ALABAMA:

35 Section 1. Section 20-2-23, Code of Alabama 1975, is
36 amended to read as follows:

37 "§20-2-23

38 (a) The Legislature finds the following:

39 (1) New synthetic substances are being created which
40 are not controlled under the provisions of existing state law
41 but which have a potential for abuse similar to or greater
42 than that for substances controlled under existing state law.
43 These new synthetic substances are called "synthetic
44 controlled substances or synthetic controlled substance
45 analogues" and can be designed to produce a desired
46 pharmacological effect and to evade the controlling statutory
47 provisions. Synthetic controlled substances or synthetic
48 controlled substance analogues are being manufactured,
49 distributed, possessed, and used as substitutes for controlled
50 substances.

51 (2) The hazards attributable to the traffic in and use
52 of a synthetic controlled substance or synthetic controlled
53 substance analogues are increased because their unregulated
54 manufacture produces variations in purity and concentration.

55 (3) Many new synthetic substances are untested, and it
56 cannot be immediately determined whether they have useful



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57 medical or chemical purposes.

58 (4) The uncontrolled importation, manufacture,
59 distribution, possession, or use of controlled substance
60 analogues has a substantial and detrimental impact on the
61 health and safety of the people of this state.

62 (5) Synthetic controlled substances or synthetic
63 controlled substance analogues can be created more rapidly
64 than they can be identified and controlled by action of the
65 Legislature. There is a need for a speedy determination of
66 their proper classification under existing law. It is
67 therefore necessary to identify and classify new substances
68 that have a potential for abuse, so that they can be
69 controlled in the same manner as other substances controlled
70 under existing state law.

71 (b) The controlled substances listed in this section
72 are included in Schedule I:

73 (1) Any of the following opiates, including their
74 isomers, esters, ethers, salts, and salts of isomers, esters,
75 and ethers, unless specifically excepted, whenever the
76 existence of these isomers, esters, ethers, and salts is
77 possible within the specific chemical designation:

- 78 a. Acetylmethadol;
- 79 b. Allylprodine;
- 80 c. Alphacetylmethadol;
- 81 d. Alphameprodine;
- 82 e. Alphamethadol;
- 83 f. Benzethidine;
- 84 g. Betacetylmethadol;



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85 h. Betameprodine;
86 i. Betamethadol;
87 j. Betaprodine;
88 k. Clonitazene;
89 l. Dextromoramide;
90 m. Dextrorphan;
91 n. Diampromide;
92 o. Diethylthiambutene;
93 p. Dimenoxadol;
94 q. Dimepheptanol;
95 r. Dimethylthiambutene;
96 s. Dioxaphetyl butyrate;
97 t. Dipipanone;
98 u. Ethylmethylthiambutene;
99 v. Etonitazene;
100 w. Etoxeridine;
101 x. Furethidine;
102 y. Hydroxypethidine;
103 z. Ketobemidone;
104 aa. Levomoramide;
105 bb. Levophenacylmorphane;
106 cc. Morpheridine;
107 dd. Noracymethadol;
108 ee. Norlevorphanol;
109 ff. Normethadone;
110 gg. Norpipanone;
111 hh. Phenadoxone;
112 ii. Phenampromide;



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- 113 jj. Phenomorphan;
- 114 kk. Phenoperidine;
- 115 ll. Piritramide;
- 116 mm. Proheptazine;
- 117 nn. Properidine;
- 118 oo. Racemoramide;
- 119 pp. Trimeperidine.

120 (2) Any of the following opium derivatives, their
121 salts, isomers, and salts of isomers, unless specifically
122 excepted, whenever the existence of these salts, isomers, and
123 salts of isomers is possible within the specific chemical
124 designation:

- 125 a. Acetorphine;
- 126 b. Acetyldihydrocodeine;
- 127 c. Benzylmorphine;
- 128 d. Codeine methylbromide;
- 129 e. Codeine-N-Oxide;
- 130 f. Cyprenorphine;
- 131 g. Desomorphine;
- 132 h. Dihydromorphine;
- 133 i. Etorphine;
- 134 j. Heroin;
- 135 k. Hydromorphenol;
- 136 l. Methyldesorphine;
- 137 m. Methyldihydromorphine;
- 138 n. Morphine methylbromide;
- 139 o. Morphine methylsulfonate;
- 140 p. Morphine-N-Oxide;



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- 141 q. Myrophine;
- 142 r. Nicocodeine;
- 143 s. Nicomorphine;
- 144 t. Normorphine;
- 145 u. Pholcodine;
- 146 v. Thebacon.

147 (3) Any material, compound, mixture, l or preparation
148 which contains any quantity of the following hallucinogenic
149 substances, their salts, isomers, l and salts of isomers, unless
150 specifically excepted, whenever the existence of these salts,
151 isomers, l and salts of isomers is possible within the specific
152 chemical designation:

- 153 a. 3,4-methylenedioxy amphetamine;
- 154 b. 5-methoxy-3,4-methylenedioxy amphetamine;
- 155 c. 3,4,5-trimethoxy amphetamine;
- 156 d. Bufotenine;
- 157 e. Diethyltryptamine;
- 158 f. Dimethyltryptamine;
- 159 g. 4-methyl-2,5-dimethoxy amphetamine;
- 160 h. Ibogaine;
- 161 i. Lysergic acid diethylamide;
- 162 j. ~~Marihuana~~Marijuana;
- 163 k. Mescaline;
- 164 l. Peyote;
- 165 m. N-ethyl-3-piperidyl benzilate;
- 166 n. N-methyl-3-piperidyl benzilate;
- 167 o. Psilocybin;
- 168 p. Psilocyn;



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169 q. Tetrahydrocannabinols, except for
170 ~~tetrahydrocannabinols~~ non-psychoactive cannabinoids derived
171 from or found in hemp, as defined in Section 2-8-381.

172 (4)a. A synthetic controlled substance that is any
173 material, mixture, or preparation that contains any quantity
174 of the following chemical compounds, their salts, isomers, and
175 salts of isomers, unless specifically excepted, whenever the
176 existence of these salts, isomers, and salts of isomers is
177 possible within the specific chemical designation or compound:

178 1. 3,4-Methylenedioxymethcathinone (Methylone), some
179 trade or other names: 3,4-methylenedioxy-N-methylcathinone.

180 2. 3,4-Methylenedioxypropylone, some other trade
181 names: (MDPV).

182 3. 4-Methylmethcathinone (Mephedrone), some trade or
183 other names: 4-methylephedrone.

184 4. 4-Methoxymethcathinone (Methedrone), some trade or
185 other names: bk-PMMA.

186 5. 3-Fluoromethcathinone, some trade or other names:
187 3-FMC.

188 6. 4-Fluoromethcathinone (Flephedrone), some trade or
189 other names: 4-FMC.

190 7.

191 1-[(5-fluoropentyl)-1H-indol-3-yl]-(2-iodophenyl)methanone,
192 some trade or other names: AM-694.

193 8.

194 1-[(5-fluoropentyl)-1H-indol-3-yl]-(naphthalen-1-yl)methanone,
195 some trade or other names: AM-2201.

196 9. (6aR, 10aR)-9-(hydroxymethyl)-6,



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197 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[
198 c]chromen-1-ol, some trade or other names: HU-210.

199 10.

200 (6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-y
201 l)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol, some trade or
202 other names: HU-211, Dexanabinol.

203 11. 1-Pentyl-2-methyl-3-(1-naphthoyl)indole, some trade
204 or other names: JWH-007.

205 12.

206 (2-Methyl-1-propyl-1H-indol-3-yl)-1-naphthalenylmethanone,
207 some trade or other names: JWH-015.

208 13. Naphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
209 trade or other names: JWH-018.

210 14. 1-Hexyl-3-(naphthalen-1-oyl)indole, some trade or
211 other names: JWH-019.

212 15. Naphthalen-1-yl-(butylindol-3-yl)methanone, some
213 trade or other names: JWH-073.

214 16.

215 4-Methoxynaphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
216 trade or other names: JWH-081.

217 17.

218 4-Methoxynaphthalen-1-yl-(1-pentyl-2-methylindol-3-yl)
219 methanone, some trade or other names: JWH-098.

220 18.

221 4-Methylnaphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
222 trade or other names: JWH-122.

223 19.

224 (1-(2-Morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-ylmethanone



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225 , some trade or other names: JWH-200.

226 20. 2-(2-Chlorophenyl)-1-(1-pentylindol-3-yl)ethanone,
227 some trade or other names: JWH-203.

228 21.

229 4-Ethyl-naphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
230 trade or other names: JWH-210.

231 22. 2-(2-Methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone,
232 some trade or other names: JWH-250.

233 23.

234 5-(2-fluorophenyl)-1-pentylpyrrol-3-yl-naphthalen-1-ylmethano
235 ne, some trade or other names: JWH-307.

236 24. 1-Pentyl-3-(4-Chloro-1-naphthoyl)indole, some trade
237 or other names: JWH-398.

238 25.

239 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol
240 (Cannabicyclohexanol), some trade or other names: CP 47, 497,
241 and homologues.

242 26.

243 2-(2-Methoxyphenyl)-1-[1-(2-cyclohexylethyl)indol-3-yl]ethanon
244 e, some trade or other names: RCS-8, SR-18.

245 27.

246 2-(4-Methoxyphenyl)-1-(1-pentyl-indol-3-yl)methanone, some
247 trade or other names: RCS-4.

248 28.

249 (R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1
250 ,2,3-de]-1,4-benzoxazin-6-yl]-1-naphthalenylmethanone, some
251 trade or other names: WIN 55,212-2.

252 29.



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253 (4-Methoxyphenyl)-[2-methyl-1-(2-morpholin-4-ylethyl)indol-3-y
254 l]methanone, some trade or other names: WIN 48,098,
255 Pravadoline.

256 b. In addition to any material, mixture, or preparation
257 that contains any quantity of the chemical compounds listed in
258 paragraph a., a synthetic controlled substance also includes
259 the following chemical compounds, their salts, isomers, and
260 salts of isomers, unless specifically excepted, whenever the
261 existence of these salts, isomers, and salts of isomers is
262 possible within the specific chemical designation or compound:

263 1.

264 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole,
265 some trade or other names: (AM-2233).

266 2. 1-Pentyl-3-(1-adamantoyl)indole, some trade or other
267 names: (AB001).

268 3.

269 [1-[(1-methyl-2-piperidiny)methyl]-1H-indol-3-yl]-1-naphthale
270 nyl-methanone, some trade or other names: (AM1220).

271 4.

272 1-(5-Fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole,
273 some trade or other names: (XLR11).

274 5. 1-Pentyl-3-(2,2,3,3-tetramethylcyclopropoyl)indole,
275 some trade or other names: (UR-144).

276 6.

277 6-Methyl-2[(4-methylphenyl)amino]-4H-3,1-benzoxazin-4-one,
278 some trade or other names: (URB 754).

279 7. [1,1'-biphenyl]-3-yl-carbamic acid, cyclohexyl
280 ester, some trade or other names: (URB 602).



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- 281 8.
282 (3'-(Aminocarbonyl)[1,1'-biphenyl]-3-yl)-cyclohexylcarbamate,
283 some trade or other names: (URB597).
- 284 9. 1-(5-Fluoropentyl)-3-(4-methyl-1-naphthoyl)indole,
285 some trade or other names: (MAM2201).
- 286 10.
287 1-naphthalenyl[4-(pentyloxy)-naphthalenyl]methanone, some
288 trade or other names: (CB-13).
- 289 11.
290 1-(5-Chloropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole,
291 some trade or other names: (5-Chloro-UR-144).
- 292 12.
293 1-(5-Fluoropentyl)-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indole-3-
294 carboxamide, some trade or other names: (STS-135).
- 295 13.
296 1[(N-Methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl)indole,
297 some trade or other names: (AM1248).
- 298 14. N-Adamantyl-1-pentyl-1H-indole-3-carboxamide, some
299 trade or other names: (SDB-001, 2NE1).
- 300 15.
301 1-Pentyl-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indazole-3-carboxam
302 ide, some trade or other names: (AKB48, APINACA).
- 303 16. 3-Naphthoylindole.
- 304 17.
305 1-[2-(4-Morpholinyl)ethyl]-3-(2,2,3,3-tetramethylcyclopropyl) i
306 ndole, some trade or other names: (A 796,260).
- 307 18.
308 1-[(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetram



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309 ethylcyclopropyl)methanone, some trade or other names: (A
310 834,735).

311 19. 1-(Pent-4-en-1-yl)-3-(4-methyl-1-naphthoyl)indole,
312 some trade or other names: (JWH-122 4-pentenyl analog).

313 20.

314 N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-fluorophenyl)m
315 ethyl]-1H-indazole-3-carboxamide some trade or other names:
316 (AB-FUBINACA).

317 21.

318 [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcycloprop
319 yl)methanone, some trade or other names: (5-Bromo-UR-144)

320 22.

321 5-(1,1-Dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexylphenol,
322 some trade or other names: (CP-47,497 C8 homolog).

323 23.

324 1-(5-Fluoropentyl)-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indazole-
325 3-carboxamide, some trade or other names: (5F-AKB48,
326 5F-APINACA).

327 24. 1-(penta-4-ene)-3-(1-naphthoyl)indole, some trade
328 or other names: (JWH-022).

329 25. 1-(5-Chloropentyl)-3-(1-naphthoyl)indole, some
330 trade or other names: (Chloro-AM-2201, JWH-018
331 N-5-chloropentyl analog).

332 26. 1-(5-Hydroxypentyl)-3-(1-naphthoyl)indole, some
333 trade or other names: (Hydroxy-AM-2201).

334 27.

335 N-[(2E)-3-(2-Methoxyethyl)4,5-dimethyl-1,3-thiazole-2(3H)-ylid
336 ene]-2,2,3,3-tetramethylcyclopropane carboxamide, some trade



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337 or other names: (A 836,339).

338 28. 1-Pentyl-3-(2-iodobenzoyl)indole, some trade or
339 other names: (AM 679).

340 29. 1-Pentyl-3-(2-methylphenacetyl)indole, some trade
341 or other names: (JWH-251).

342 30. 1-pentyl-1H-indole-3-carboxylic acid 8-quinolinyl
343 ester, some trade or other names: (PB-22, QUPIC).

344 31. 1-(5-fluoropentyl)-1H-indole-3-carboxylic acid
345 8-quinolinyl ester, some trade or other names: (5F-PB-22).

346 32.

347 1-pentyl-N-(naphthalen-1-yl)-1H-indole-3-carboxamide, some
348 trade or other names: (MN-24, NNE1).

349 33. 1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid
350 8-quinolinyl ester, some trade or other names: (BB-22,
351 QUCHIC).

352 34.

353 N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-pentyl-1H-indazole-
354 -3-carboxamide, some trade or other names: (AB-PINACA).

355 35.

356 7-methoxy-1-(2-morpholinoethyl)-N-((1S,2S,4R)-1,3,3-trimethylb
357 icyclo[2.2.1]heptan-2-yl)-1H-indole-3-carboxamide, some trade
358 or other names: (MN-25).

359 36. ADB-PINACA.

360 37. FUB-AKB-48.

361 38. FUB-PB-22.

362 39. Heptyl-UR144.

363 40. THJ-018.

364 41. THJ-2201.



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365 42. 1-heptyl-3-(1-naphthoyl)indole), some trade or other
366 names: (JWH-20).

367 43. Napthalen-1-yl-(1-propyl-1H-indol-3-yl)methanone,
368 some trade or other names: (JWH-072).

369 44. (6aR,10aR)-3-(1, 1-Dimethylbutyl)-6a, 7, 10,
370 10a-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran, some
371 trade or other names: (JWH-133).

372 45. 3-(naphthalen-1-ylmethyl)-1-pentyl-1H-indole, some
373 trade or other names: (JWH-175).

374 46. 1-pentyl-3-(4-methoxyphenylacetyl)indole, some
375 trade or other names: (JWH-201).

376 47. 1-pentyl-3-(3-methoxyphenylacetyl)indole, some
377 trade or other names: (JWH 302).

378 48.

379 [(1R,2R,5R)-2-[2,6-dimethoxy-4-(2-methyloctan-2-yl)phenyl]-7,7-
380 -dimethyl-4-bicyclo[3.1.1]hept-3-enyl]methanol, some trade or
381 other names: (HU-308).

382 49.

383 3-hydroxy-2-[(1R,6R)-3-methyl-6-(1-methylethenyl)-2-cyclohexen-
384 -1-yl]-5-pentyl-2,5-cyclohexadiene-1,4-dione, some trade or
385 other names: (HU-331).

386 50.

387 N-cyclopropyl-11-(3-hydroxy-5-pentylphenoxy)-undecanamide,
388 some trade or other names: (CB-25).

389 51.

390 N-cyclopropyl-11-(2-hexyl-5-hydroxyphenoxy)-undecanamide, some
391 trade or other names: (CB-52).

392 52.



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393 2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-5-(2-me
394 thyloctan-2-yl)phenol, some trade or other names:
395 (CB-55,940) (CB-55).

396 53. 4-Methylethylcathinone, some trade or other names:
397 (4-MEC, 4-Methylethcathinone).

398 54. 4'-Methyl-alpha-pyrrolidinopropiophenone, some
399 trade or other names: (MPPP, ZZ-1).

400 55.
401 (RS)-1-naphthalen-2-yl-2-pyrrolidin-1-ylpentan-1-one, some
402 trade or other names: (Naphyrone).

403 56. alpha,alpha-Diphenyl-2-piperidinemethanol, some
404 trade or other names: (Pipradrol, Meratran).

405 57.
406 (RS)-1-(4-methylphenyl)-2-(1-pyrrolidinyl)pentan-1-one, some
407 trade or other names: (Pyrovalerone).

408 58. 3,4-Dimethylmethcathinone, some trade or other
409 names: (3,4-DMMC).

410 59. 4-Fluoroamphetamine, some trade or other names:
411 (4-FA).

412 60. 4-Fluoromethamphetamine, some trade or other names:
413 (4-FMA).

414 61. Butylone, some trade or other names: (bk-MBDB).

415 62. alpha-Pyrrolidinopentiophenone, some trade or other
416 names: (alpha-PVP).

417 63. beta-keto-Dimethylbenzodioxolylbutanamine, some
418 trade or other names: (bk-DMBDB).

419 64. 2-(methylamino)-1-phenylbutan-1-one, some trade or
420 other names: (Buphedrone).



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- 421 65. (RS)-2-ethylamino-1-phenyl-propan-1-one, some trade
422 or other names: (N-Ethylcathinone).
- 423 66. 2-Fluoroamphetamine, some trade or other names:
424 (2-FA).
- 425 67. Methoxetamine, some trade or other names: (MXE).
- 426 68. 2-Methylamino-1-phenylpentan-1-one, some trade or
427 other names: (Pentedrone).
- 428 69. 3,4-Methylenedioxcathinone, some trade or other
429 names: (MDC).
- 430 70. 2-Fluoromethamphetamine, some trade or other names:
431 (2-FMA).
- 432 71. 4-methylmethamphetamine, some trade or other names:
433 (4-MMA).
- 434 72. 4-Fluoroisocathinone, some trade or other names:
435 (4-FIC).
- 436 73. 3-Fluoromethamphetamine, some trade or other names:
437 (3-FMA).
- 438 74. Methiopropamine, some trade or other names: (MPA).
- 439 75. alpha-Pyrrolidinobutiophenone, some trade or other
440 names: (alpha-PBP).
- 441 76. 4-Methoxy-N-methylcathinone, some trade or other
442 names: (Methedrone, bk-PMMA).
- 443 77. alpha-Pyrrolidinopropiophenone, some trade or other
444 names: (alpha-PPP).
- 445 78. (RS)-2-benzhydrylpiperidine, some trade or other
446 names: (Desoxypipradrol).
- 447 79. 3,4-Methylenedioxyethylcathinone, some trade or
448 other names: (MDEC).



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- 449 80. 3,4-Methylenedioxy-alpha-pyrrolidinobutiophenone,
450 some trade or other names: (MDPBP).
- 451 81.
452 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one
453 (Pentylone, bk-MBDP).
- 454 82. 3-Fluoroamphetamine, some trade or other names:
455 (3-FA).
- 456 83. 3-Fluoromethcathinone, some trade or other names:
457 (3-FMC).
- 458 84. 2-Fluoromethcathinone, some trade or other names:
459 (2-FMC).
- 460 85.
461 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)propan-1-one
462 (bk-MDDMA).
- 463 86. N,N-Diethylcathinone, some trade or other names:
464 (Amfepramone, DEC).
- 465 87. 1,3-Dimethylamylamine, some trade or other names:
466 (DMAA).
- 467 88. N, N-Dimethylcathinone, some trade or other names:
468 (DMC).
- 469 89. N-Ethyl-3,4-methylenedioxycathinone, some trade or
470 other names: (bk-MDEA).
- 471 90. N-Ethylamphetamine, some trade or other names:
472 (EMA).
- 473 91. N-Ethylcathinone, some trade or other names: (EC).
- 474 92. 2-Ethylethcathinone, some trade or other names:
475 (2-EEC).
- 476 93. 4-Ethyl-N-ethylcathinone, some trade or other



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477 names: (4-EEC) .
478 94.
479 2-(5-Methoxy-1-benzofuran-3-yl)-N,N-dimethylethanamine, some
480 trade or other names: (Dimembfe) .
481 95. 2-(5-Methoxy-1-benzofuran-3-yl)N-ethylethamine.
482 96. 4-Methoxymethamphetamine, some trade or other
483 names: (PMMA) .
484 97. 4-Methoxy-N-ethylamphetamine, some trade or other
485 names: (PMEA) .
486 98. 4-Methoxy-N-ethylcathinone, some trade or other
487 names: (ETHEDRONE) .
488 99. 3-Methylmethcathinone, some trade or other names:
489 (3-MMC) .
490 100. 4-Methyl-alpha-pyrrolidinobutiophenone, some trade
491 or other names: (MPBP) .
492 101. 2-Methylethcathinone, some trade or other names:
493 (2-MEC) .
494 102. 3-Methylethcathinone, some trade or other names:
495 (3-MEC) .
496 103. 2-Ethylethcathinone, some trade or other names:
497 (2-EEC) .
498 104. 3-Ethylethcathinone, some trade or other names:
499 (3-EEC) .
500 105. 3-Ethylmethcathinone, some trade or other names:
501 (3-EMC) .
502 106.
503 3',4'-Methylenedioxy-alpha-pyrrolidinopropiophenone, some
504 trade or other names: (MDPPP) .



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- 505 107. alpha-Pyrrolidinopentiothiophenone, some trade or
506 other names: (alpha-PVT).
- 507 108. 3-Methoxymethcathinone, some trade or other names:
508 (3-MeOMC).
- 509 109. N-Methyl-1,3-benzodioxolylbutanamine, some trade
510 or other names: (MBDB).
- 511 110. Ethcathinone, some trade or other names:
512 (ETHYLPROPION, ETH-CAT).
- 513 111. Ethylone (3,4-methylenedioxy-N-ethylcathinone).
- 514 112. N-N-Diethyl-3,4-methylenedioxy-cathinone.
- 515 113. 3,4-methylenedioxy-propiofenone.
- 516 114. 2-Bromo-3,4-methylenedioxypropiofenone.
- 517 115. 3,4-methylenedioxy-propiofenone-2-oxime.
- 518 116. N-Acetyl-3,4-methylenedioxy-cathinone.
- 519 117. N-Acetyl-N-Methyl-3,4-methylenedioxy-cathinone.
- 520 118. N-Acetyl-N-Ethyl-3,4-methylenedioxy-cathinone.
- 521 119. 4-Bromomethcathinone.
- 522 120. 3-Bromomethcathinone.
- 523 121. Eutylone (beta-Keto-Ethylbenzodioxolylbutanamine).
- 524 122. 4'-Methoxy-alpha-pyrrolidinopropiofenone, some
525 trade or other names: (MOPPP).
- 526 123. 4'-Methyl-alpha-pyrrolidinohexiofenone, some
527 trade or other names: (MPHP).
- 528 124. Benocyclidine (BCP) or
529 Benzothiophenylcyclohexylpiperidine, some trade or other
530 names: (BTCP).
- 531 125. 4-Fluoro-(methylamino)butyrophenone, some trade or
532 other names: (F-MABP).



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- 533 126. 3-Methyl-4-Methoxymethacathinone, some trade or
534 other names: (3-Me-4-MeO-MCAT).
- 535 127. 4-Methyl-(ethylamino)-butryophenone, some trade or
536 other names: (Me-EABP).
- 537 128. 4-Ethyl-methcathinone, some trade or other names:
538 (4-EMC).
- 539 129. 4-methoxy-N-ethylcathinone
540 (bk-PMC;p-methox-ethcathinone).
- 541 130. 4'-Methoxy-alpha-pyrroldino-propiofenone (MeOPPP;
542 4'-MeO-PPP).
- 543 131. 3-Fluorocathinone (3-FC).
- 544 132. 4-Fluorocathinone (4-FC).
- 545 133. 4-methyl-buphedrone (4-MeMABP; 4MeBP; BZ-6378).
- 546 134. 3,4-Methylenedioxy-N-benzylcathinone, some trade
547 or other names: (BMDP).
- 548 135. N-Benzyl-butylone, some trade or other names:
549 (BMDB).
- 550 136. N-Hydroxy-3,4-methylenedioxymethcathinone.
- 551 137. N-ethylbuphedrone, some trade or other names:
552 (NEB).
- 553 138. 4-Fluorobuphedrone, some trade or other names:
554 (4-FBP).
- 555 139. 4-Methoxy-pyrrolidinbutrophenone (4-MeO-PBP).
- 556 140. 4-Ethyl-pyrrolidinobutrophenone, some trade or
557 other names: (4-Et-PBP).
- 558 141. 5-(2-aminopropyl)indole, some trade or other
559 names: (5-IT).
- 560 142. 1-phenyl-2-(piperidin-1-yl)butan-1-one.



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561 143. 2,4,5-Trimethyl-methacathinone, some trade or
562 other names: (2,4,5-TMMC).

563 144. alpha-pyrrolidino-heptiophenone, some trade or
564 other names: (alpha-PHpP).

565 145. 4-Methylamphetamine (4-MA: pTAP; PAL-313; 4-MeA;
566 PmeA).

567 146. N-Ethyl-methamphetamine.

568 147. 4-(2-Aminopropyl)benzofuran, some trade or other
569 names: (4-APB).

570 148. 5-(2-Aminopropyl)-2,3-dihydro-1H-indene (5-APDI;
571 IAP; AIP; indanylaminoporpane).

572 149. 6,7-Methylenedioxy-2-aminotetralin, some trade or
573 other names: (MDAT).

574 150. 4-Methylthioamphetamine (4-MTA; P1882).

575 151. 4-Chloroamphetamine (p-chloro-amphetamine).

576 152. 2,4,6-Trimethoxyamphetamine, some trade or other
577 names: (TMA-6).

578 153. 2,4,5-Trimethoxyamphetamine, some trade or other
579 names: (TMA-2).

580 154. 2,5-Dimethylamphetamine, some trade or other
581 names: (2,5-DMA).

582 155. 3,4-Dimethylamphetamine, some trade or other
583 names: (3,4-DMA).

584 156. N-propylamphetamine.

585 157. 4-Hydroxyamphetamine.

586 158. 3-Hydroxyamphetamine.

587 159. Methylenedioxydimethylamphetamine, some trade or
588 other names: (MDDM).



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- 589 160. 2-Aminoindane, some trade or other names: (2-AI).
- 590 161. 5,6-Methylenedioxy-N-methyl-aminoindane, some
591 trade or other names: (MDMAI).
- 592 162. 2C-T-21.
- 593 163. 2C-B-Fly.
- 594 164. 3,4-dimethyl-2,5-dimethoxyphenethylamine (2C-G).
- 595 165. 25D-NBOMe.
- 596 166. 25G-NBOMe.
- 597 167. 25N-NBOMe.
- 598 168. Bromo-benzylidifuranyl-isopropylamine, some trade
599 or other names: (Bromo Dragon Fly).
- 600 169. 3C-B fly.
- 601 170. 2,5-Dimethoxy-4-ethylthioamphetamine, some trade
602 or other names: (Aleph-2).
- 603 171. 1-[(4-ethoxy-2,5-dimethoxy)phenyl]propan-2-amine,
604 some trade or other names: (MEM).
- 605 172.
- 606 1-[2,5-dimethoxy-4-(propylthio)phenyl]propan-2-amine, some
607 trade or other names: (Aleph-7).
- 608 173. N-benzyl-2-phenylethylamine.
- 609 174. N,N-dimethyl-2-phenylethylamine.
- 610 175. 6-chloro-2-aminotetralin, some trade or other
611 names: (6-CAT).
- 612 176. 2-phenylpropan-1-amine, some trade or other names:
613 (B-Me-PEA).
- 614 177. 2-Phenethylamine, some trade or other names:
615 (2-PEA).
- 616 178. 1-methylamino-1-(3,4-methylenedioxyphenyl)propane,



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617 some trade or other names: (M-ALPHA).
618 179. Camfetamine.
619 180. Methoxyphenamine.
620 181. 4-methylaminorex, some trade or other names:
621 (4-MAR; 4-MAX; U4Euh; Euphoria; Ice).
622 182. (1-thiophen-2-yl)propan-2-amine
623 (Thienoamphetamine).
624 183. Dimethocaine.
625 184. 4-Fluoroephedrine.
626 185. 4-methyaminorex (p-methyl derivative).
627 186.
628 1-[(N-methylpiperidin-2-yl)methyl]-2-methyl-3-(naphthalen-1-oy
629 l)-6-nitroindole (AM1221).
630 187. (1-butyl-1H-indol-3-yl)(4-methoxyphenyl)-methanone
631 (RCS-4 (C4) homolog).
632 188. 5-[3-(1-naphthoyl)-1H-indole-1-yl]pentanenitrile,
633 some trade or other names: (AM2232).
634 189. 1-(Pentyl)-3-(4-bromo-1-naphthoyl)-indole, some
635 trade or other names: (JWH-387).
636 190. 1-(Pentyl)-3-(4-fluoro-1-naphthoyl)-indole, some
637 trade or other names: (JWH-412).
638 191. 1-(5-chloropentyl)-3-(2-iodobenzoyl)indole, some
639 trade or other names: (AM694 Derivative).
640 192.
641 (2-iodo-5-nitrophenyl)-[1-[(1-methylpiperidin-2-yl)methyl]1H-i
642 ndol-3-yl]-methanone, some trade or other names: (AM1241).
643 193. 1-Pentyl-3-[1-(4-propyl)naphthoyl]indole, some
644 trade or other names: (JWH-182).



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- 645 194. JWH-081 2-methoxynaphthyl isomer, some trade or
646 other names: (JWH-267).
- 647 195.
648 (3-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone, some trade
649 or other names: (RCS-4 3-methoxy isomer).
- 650 196.
651 [1-(5-fluoropentyl)-1H-indol-3-yl](4-ethyl-1-naphthalenyl)-met
652 hanone (EAM-2201).
- 653 197. ADB-FUBINACA.
- 654 198. ADBICA.
- 655 199. AM-279.
- 656 200. JWH-370.
- 657 201. NNE-1.
- 658 202. MAM-2201 chloropentyl derivative.
- 659 203. 1-(5-fluoropentyl)-3-(2-methyl-benzoyl)indole.
- 660 204. 1-(5-fluoropentyl)-3-(2-ethylbenzoyl)indole.
- 661 205. AB-005.
- 662 206. AB-005 Azepane isomer.
- 663 207.
664 4-hydroxy-3,3,4-trimethyl-1-(1-pentyl-1H-indol-3-yl)pentan-1-o
665 ne (4-HTMPIPO).
- 666 208. UR-12.
- 667 209. 5-Fluoro-ADBICA.
- 668 210. BAY-38-7271; KN 38-7271.
- 669 211. JTE-907.
- 670 212. Org 27569.
- 671 213. Org 27759.
- 672 214. Org 29647.



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- 673 215. LY 2183240.
- 674 216. JTE 7-31.
- 675 217. URB 937.
- 676 218. 3-methoxy-eticyclidine, some trade or other names:
677 (3-MeO-PCE).
- 678 219. 1-Phenylcyclohexanamine, some trade or other
679 names: (PCA).
- 680 220. 4-Methyl-phencyclidine, some trade or other names:
681 (4-Me-PCP).
- 682 221. 4-Methoxy-eticyclidine, some trade or other names:
683 (4-MeO-PCE).
- 684 222. 4-Methoxyphencyclidine, some trade or other names:
685 (Methoxydine; 4MeO-PCP).
- 686 223. 3-Methoxyphencyclidine, some trade or other names:
687 (3-MeO-PCP).
- 688 224. 1-phenyl-N-propylcyclohexanamine, some trade or
689 other names: (PCPr).
- 690 225. N-(2-methoxyethyl)-1-phenylcyclohexanamine, some
691 trade or other names: (PCMEA).
- 692 226. N-(2-ethoxyethyl)-1-phenylcyclohexanamine, some
693 trade or other names: (PCEEA).
- 694 227. N-(3-methoxypropyl)-1-phenylcyclohexanamine, some
695 trade or other names: (PCMPA).
- 696 228. 3-Hydroxy-phencyclidine, some trade or other
697 names: (3-OH-PCP).
- 698 229. Methoxyketamine, some trade or other names:
699 (2-MeO-2-deschloro-ketamine).
- 700 230. Tiletamine, some trade or other names: (TCE).



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701 231. N-ethylnorketamine.

702 232. N-Methyltryptamine, some trade or other names:
703 (NMT).

704 233. N-Methyl-N-isopropyltryptamine, some trade or
705 other names: (MiPT; MIPT).

706 234. 4-hydroxy-N,N-methylisopropyltryptamine, some
707 trade or other names: (4-OH-MiPT).

708 235. 4-Acetoxy-N,N-diisopropyl-tryptamine (4-AcO-DiPT:
709 4-AcO-DIPT; 4-Acetoxy-MiPT).

710 236. 4-Methoxy-N,N-dimethyltryptamine, some trade or
711 other names: (4-MeO-DMT).

712 237. 5-Hydroxytryptamine, some trade or other names:
713 (5-HT).

714 238. 5-acetoxy-N,N-dimethyltryptamine, some trade or
715 other names: (5-AcO-DMT).

716 239. 5-Methoxy-N,N-dipropyltryptamine, some trade or
717 other names: (5-MeO-DPT).

718 240. d-Lysergic acid amide, some trade or other names:
719 (LSA; ergine).

720 241. 2,5-dimethoxy-4-chloroamphetamine, some trade or
721 other names: (DOC).

722 242.

723 N-(2-Methoxybenzyl)-4-iodo-2,5-dimethoxyphenethylamine, some
724 trade or other names: (25I-NBOMe).

725 243. 4-Ethyl-2,5-dimethoxyphenethylamine, some trade or
726 other names: (2C-E).

727 244. 2,5-Dimethoxy-4-iodophenethylamine, some trade or
728 other names: (2C-I).



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729 245. 6-(2-Aminopropyl)-2,3-dihydrobenzofuran, some
730 trade or other names: (6-APDB).

731 246. 6-(2-Aminopropyl)benzofuran, some trade or other
732 names: (6-APB).

733 247. 5-(2-Aminopropyl)-2,3-dihydrobenzofuran, some
734 trade or other names: (5-APDB).

735 248. 5-(2-Aminopropyl)benzofuran, some trade or other
736 names: (5-APB).

737 249. 2,5-Dimethoxy-4-(n)-propylthiophenethylamine, some
738 trade or other names: (2C-T-7).

739 250. 2,5-Dimethoxy-4-(n)-propylphenethylamine, some
740 trade or other names: (2C-P).

741 251. 2,5-Dimethoxy-4-bromoamphetamine, some trade or
742 other names: (DOB).

743 252. 2,5-Dimethoxy-4-bromobenzylpiperazine, some trade
744 or other names: (2C-B-BZP).

745 253. 2,5-Dimethoxy-4-bromophenethylamine, some trade or
746 other names: (2C-B).

747 254. 2,5-Dimethoxy-4-chlorophenethylamine, some trade
748 or other names: (2C-C).

749 255. 2,5-Dimethoxy-(4-ethylthio)phenethylamine, some
750 trade or other names: (2C-T-2).

751 256. 2,5-Dimethoxy-4-iodoamphetamine, some trade or
752 other names: (DOI).

753 257. 2,5-Dimethoxy-4-methylamphetamine, some trade or
754 other names: (DOM).

755 258. 2,5-Dimethoxyphenethylamine, some trade or other
756 names: (2C-H).



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757 259.
758 2-(2,5-Dimethoxyphenyl-4-bromo)-N-(2-methoxybenzyl)ethanamine,
759 some trade or other names: (25B-NBOMe).
760 260.
761 2-(2,5-Dimethoxyphenyl-4-chloro)-N-(2-methoxybenzyl)ethanamine
762 , some trade or other names: (25C-NBOMe).
763 261.
764 2-(2,5-Dimethoxyphenyl-4-ethyl)-N-(2-methoxybenzyl)ethanamine,
765 some trade or other names: (25E-NBOMe).
766 262. 2-Ethylmethcathinone, some trade or other names:
767 (2-EMC).
768 263.
769 2-(2,5-Dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, some
770 trade or other names: (25H-NBOMe).
771 264. BZP (Benzylpiperazine).
772 265. para-Fluorophenylpiperazine.
773 266. 1-(4-Methylphenyl)piperazine.
774 267. meta-Cholorophenylpiperazine.
775 268. para-Methoxyphenylpiperazine.
776 269. DBZP (1,4-dibenzylpiperazine).
777 270. TFMPP (3-Trifluoromethylphenylpiperazine).
778 271. 2C-T-4
779 (2,5-Dimethoxy-4-isopropylthiophenethylamine).
780 272. 2C-T (2,5-Dimethoxy-4-methylthiophenethylamine).
781 273. 2C-D (2-(2,5-Dimethoxy-4-methylphenyl)ethanamine).
782 274. 2C-N 2,5-Dimethoxy-4-nitrophenethylamine.
783 275. 5-methoxy-N,N-diallyltryptamine, some trade or
784 other names: (5-MeO-DALT).



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785 276. 5-Methoxy-N,N-Diisopropyltryptamine, some trade or
786 other names: (5-MeO-DIPT).

787 277. 5-Methoxy-alpha-methyltryptamine, some trade or
788 other names: (5-MeO-AMT).

789 278. 4-Acetoxy-N,N-dimethyltryptamine, some trade or
790 other names: (4-AcO-DMT).

791 279. 4-Hydroxy-N,N-diethyltryptamine, some trade or
792 other names: (4-HO-DET).

793 280. 4-Hydroxy-N,N-diisopropyltryptamine, some trade or
794 other names: (4-HO-DIPT).

795 281. 4-Hydroxy-N-methyl-N-ethyltryptamine, some trade
796 or other names: (4-OH-MET).

797 282. 5-Methoxy-N,N-diethyltryptamine, some trade or
798 other names: (5-MeO-DET).

799 283. 5-Methoxy-N-methyl-N-isopropyltryptamine, some
800 trade or other names: (5-MeO-MIPT).

801 284. 4-Acetoxy-N,N-diethyltryptamine, some trade or
802 other names: (4-AcO-DET).

803 285. 4-Acetoxy-N-methyl-N-isopropyltryptamine, some
804 trade or other names: (4-AcO-MIPT).

805 286. N,N-Dipropyltryptamine, some trade or other names:
806 (DPT).

807 287. N,N-Diisopropyltryptamine, some trade or other
808 names: (DIPT).

809 288. 4-Methoxy-N-methyl-N-isopropyltryptamine, some
810 trade or other names: (4-MeO-MIPT).

811 289. Tyramine (4-Hydroxyphenethylamine).

812 290. 5-Hydroxy-alpha-methyltryptamine.



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- 813 291. 5-Hydroxy-N-methyltryptamine.
814 292. 5-Methoxy-N,N-dimethyltryptamine.
815 293. 5-Methyl-N,N-dimethyltryptamine.
816 294. Diphenylprolinol, some trade or other names:
817 (D2PM; diphenyl-2-pyrrolidinemethanol).
818 295. 3,4 Dichloromethylphenidate, some trade or other
819 names: (3,4-CTMP).
820 296. 3-chloromethyl-phenidate, some trade or other
821 names: (3-CTMP).
822 297. 4-Methylmethylphenidate.
823 298. 4-Fluoromethyl-phenidate, some trade or other
824 names: (4-FTMP).
825 299. Ethylphenidate.
826 300. Etizolam (Etilaam, Etizola, Sedekopan, Pasaden,
827 Depas).
828 301. Phenazepam.
829 302. Pyrazolam.
830 303. CL-218,872.
831 304. Zopiclone.
832 305. Salvinorin A.
833 306. AH-7921.
834 307. O-Desmethyltramadol, some trade or other names:
835 (O-DT; ODT).
836 308. Desmorphine (Dihydrodesoxymorphine; permonid;
837 krokodil; crocodile).
838 309. Acetyl Fentanyl (desmethylfentanyl).
839 310. 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine
840 (MT-45).



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- 841 311. 1-(2-methoxyphenyl)piperazine, some trade or other
842 names: (MOPIP).
- 843 312. 1-(4-Chlorophenyl)piperazine, some trade or other
844 names: (pCPP).
- 845 313. para-Methoxyphenyl-piperazine, some trade or other
846 names: (MBZP).
- 847 314. Methylmethaqualone.
- 848 315. Etaqualone.
- 849 316. 5-Iodo-2-aminoindane, some trade or other names:
850 (5-IAI).
- 851 317. 5,6-(Methylenedioxy)-2-aminoindane, some trade or
852 other names: (5,6-MDAI).
- 853 318. 4,5-(Methylenedioxy)-2-aminoindane, some trade or
854 other names: (4,5-MDAI).
- 855 319. MMAI.
- 856 320. W-15.
- 857 321. W-18.
- 858 322. Mitragynine.
- 859 323. Hydroxymitragynine.
- 860 324. Butyrfentanyl
861 (N-phenyl-N-[1-(2-phenylethyl)-4-pyridinyl]butyramide).
- 862 325. Beta-Hydroxythiofentanyl
863 (N-phenyl-N-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl}-pr
864 opanamide).
- 865 326. 4-methylphenethyl acetyl fentanyl
866 (N-phenyl-N-{1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-acetami
867 de).
- 868 327. Acrylfentanyl



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- 869 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-prop-2-enamide) .
870 328. 3-Allylfentanyl
871 (N-phenyl-N-[1-(2-phenylethyl)-(3s,4R)-3-prop-2-enyl-4-piperid
872 inyl]-propanamide) .
873 329. Benzodioxole fentanyl
874 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-benzo[d][1,3]dio
875 xole-5-carboxamide) .
876 330. Benzyl carfentanil
877 (N-phenyl-N-(1-benzyl-4-methylcarboxylate-4-piperidinyl)-propa
878 namide) .
879 331. Brifentanil
880 (N-(2-fluorophenyl)-N-((3R,4S)-1-[2-(4-ethyl-5-oxotetrazol-1-y
881 l)ethyl]-3-methyl-4-piperidinyl)-2-methoxyacetamide) .
882 332. Cyclopentylfentanyl
883 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-cyclopentanecarb
884 oxamide) .
885 333. 2,5-Dimethylfentanyl
886 (N-phenyl-N-[1-(2-phenylethyl)-2,5-dimethyl-4-piperidinyl]-pro
887 pranamide) .
888 334. 4-Fluoroisobutyryl fentanyl
889 (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-isobut
890 yramide) .
891 335. Furanyl fentanyl
892 (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]furan-2-carboxami
893 de) .
894 336. Furanylethyl fentanyl
895 (N-phenyl-N-[1-(2-furanylethyl)-4-piperidinyl]-propanamide) .
896 337. Isobutyryl fentanyl



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897 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-methylpropanami
898 de) .

899 338. Lofentanil

900 (N-phenyl-N-[1-(2-phenylethyl)-(3R,4S)-3-methyl-4-methylcarbox
901 ylate-4-piperidinyl]-propanamide) .

902 339. 4-Methoxybutyrfentanyl

903 (N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butyr
904 amide) .

905 340. 4-Methoxymethylfentanyl

906 (N-phenyl-N-[1-(2-phenylethyl)-4-methoxymethyl-4-piperidinyl]-
907 propanamide) .

908 341. Meta-fluorobutyryl fentanyl

909 (N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butyram
910 ide) .

911 342. Meta-fluorofentanyl

912 (N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propa
913 mide) .

914 343. 3-Methylbutyrfentanyl

915 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butyramide) .

916 344. N-Methylcarfentanyl

917 (N-phenyl-N-(1-methyl-4-methylcarboxylate-4-piperidinyl)-propa
918 namide) .

919 345. Methoxyacetylfentanyl

920 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-methoxyacetami
921 de) .

922 346. Mirfentanyl

923 (N-(2-pyrazinyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-2-furamid
924 e) .



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- 925 347. Ocfentanil
926 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-2-metho
927 xyacetamide).
- 928 348. Ohmefentanil
929 (N-phenyl-N-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidiny
930 l]-propanamide).
- 931 349. Ortho-fluorobutyryl fentanyl
932 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-butyram
933 ide).
- 934 350. Ortho-fluorofentanyl
935 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-propa
936 mide).
- 937 351. Para-chlorofentanyl
938 (N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propan
939 amide).
- 940 352. Para-chloroisobutyryl fentanyl
941 (N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-isobut
942 yramide).
- 943 353. 4-Fluorobutyryl fentanyl
944 (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butyra
945 mide).
- 946 354. Para-methoxyfentanyl
947 (N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propa
948 namide).
- 949 355. Para-methylfentanyl
950 (N-(4-methylphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propan
951 amide).
- 952 356. 4-Phenyl fentanyl



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953 (N-phenyl-N-[4-phenyl-1-(2-phenylethyl)-4-piperidinyl]-propana
954 mide).

955 357. Trefentanyl

956 (N-(2-fluorophenyl)-N-{1-[2-(4-ethyl-5-oxo-4,5-dihydro-1H-tetr
957 azol-1-yl)ethyl]-4-phenyl-4-piperdinyll}-propanamide).

958 358. Valeryl fentanyl

959 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide).

960 359. Alpha-Methylacetylfentanyl

961 (N-phenyl-N-[1-phenylpropan-2-yl]-4-piperidinyl]-acetamide).

962 360. Alpha-Methylbutyrfentanyl

963 (N-phenyl-N-[1-phenylpropan-2-yl]-4-piperidinyl]-butyramide).

964 361. Alpha-Methylthiofentanyl

965 (N-phenyl-N-[1-(1-thienyl-2-ylpropan-2-yl)-4-piperidinyl]-prop
966 anamide).

967 362. Beta-Hydroxy fentanyl

968 (N-phenyl-N-[1-(2-hydroxy-2-phenylethyl)-4-piperidinyl]-propan
969 amide).

970 363. Beta-Methyl fentanyl

971 (N-phenyl-N-[1-(2-phenylpropyl)-4-piperdinyll]-propanamide).

972 364. U-47700

973 (3,4-Dichloro-N-[(1R,2R)-2-(dimethylamino)cyclohexyl]-N-methyl
974 benzamide).

975 365. W-19

976 ((Z)-N-{1-[2-(4-aminophenyl)ethyl]piperidin-2-ylidene}-4-chlor
977 obenzenesulfonamide).

978 366. Flubromazolam

979 (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a]
980 [1,4]benzodiazepine).



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981 367. Tianeptine.

982 (5)a. A synthetic controlled substance analogue, being
983 a material, mixture, or preparation that contains any chemical
984 structure of which is chemically similar to the chemical
985 structure of any other controlled substance in Schedule I or
986 Schedule II or that satisfies any one of the following:

987 1. Has a stimulant, depressant, or hallucinogenic
988 effect on the central nervous system that mimics or is similar
989 to or greater than the stimulant, depressant, or
990 hallucinogenic effect on the central nervous system of a
991 controlled substance in Schedule I or Schedule II.

992 2. With respect to a particular person, if the person
993 represents or intends that the substance have a stimulant,
994 depressant, or hallucinogenic effect on the central nervous
995 system that is substantially similar to or greater than the
996 stimulant, depressant, or hallucinogenic effect on the central
997 nervous system of a controlled substance in Schedule I or
998 Schedule II and the substance is actually capable of producing
999 a stimulant, depressant, or hallucinogenic effect on the
1000 central nervous system that mimics, is similar to, or is
1001 greater than the stimulant, depressant, or hallucinogenic
1002 effect on the central nervous system of a controlled substance
1003 in Schedule I or Schedule II.

1004 3. Has been demonstrated to have binding activity at
1005 one or more cannabinoid receptors.

1006 4. Is capable of exhibiting cannabinoid-like activity.

1007 5. Any compound structurally analogous to, mimicking,
1008 or derived from 3-(1-naphthoyl)indole or



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1009 1H-indol-3-yl-(1-naphthyl)methane by substitution at the
1010 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
1011 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1012 cycloalkylmethyl, cycloalkylethyl,
1013 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1014 or 2-(4-morpholinyl)ethyl whether or not further substituted
1015 in the indole ring to any extent, whether or not substituted
1016 in the naphthyl ring to any extent.

1017 6. Any compound structurally analogous to, mimicking,
1018 or derived from 3-(1-naphthoyl)pyrrole by substitution at the
1019 nitrogen atom of the pyrrole ring by alkyl, alkyl halide, aryl
1020 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1021 cycloalkylmethyl, cycloalkylethyl,
1022 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1023 or 2-(4-morpholinyl)ethyl, whether or not further substituted
1024 in the pyrrole ring to any extent, whether or not substituted
1025 in the naphthyl ring to any extent.

1026 7. Any compound structurally analogous to, mimicking,
1027 or derived from 1-(1-naphthylmethyl)indene by substitution at
1028 the 3-position of the indene ring by alkyl, alkyl halide, aryl
1029 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1030 cycloalkylmethyl, cycloalkylethyl,
1031 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1032 or 2-(4-morpholinyl)ethyl whether or not further substituted
1033 in the indene ring to any extent, whether or not substituted
1034 in the naphthyl ring to any extent.

1035 8. Any compound structurally analogous to, mimicking,
1036 or derived from 3-phenylacetylindole by substitution at the



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1037 nitrogen atom of the indole ring with alkyl, alkyl halide,
1038 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1039 cycloalkylmethyl, cycloalkylethyl,
1040 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1041 or 2-(4-morpholinyl)ethyl, whether or not further substituted
1042 in the indole ring to any extent, whether or not substituted
1043 in the phenyl ring to any extent.

1044 9. Any compound structurally analogous to, mimicking,
1045 or derived from 2-(3-hydroxycyclohexyl)phenol by substitution
1046 at the 5-position of the phenolic ring by alkyl, alkyl halide,
1047 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1048 cycloalkylmethyl, cycloalkylethyl,
1049 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1050 or 2-(4-morpholinyl)ethyl, whether or not substituted in the
1051 cyclohexyl ring to any extent.

1052 10. Any compound structurally analogous to, mimicking,
1053 or derived from 3-(2,2,3,3-tetramethylcyclopropoyl)indole or
1054 1H-indol-3-yl-(2,2,3,3-tetramethylcyclopropoyl)methane by
1055 substitution at the nitrogen atom of the indole ring by alkyl,
1056 alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
1057 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
1058 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1059 or 2-(4-morpholinyl)ethyl whether or not further substituted
1060 in the indole ring to any extent.

1061 11. Any compound structurally analogous to, mimicking,
1062 or derived from 3-(adamant-1-oyl)indole or
1063 1H-indol-3-yl-(1-adamantyl)methane by substitution at the
1064 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl



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1065 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1066 cycloalkylmethyl, cycloalkylethyl,
1067 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1068 or 2-(4-morpholinyl)ethyl whether or not further substituted
1069 in the indole ring to any extent.

1070 12. Any compound structurally analogous to, mimicking,
1071 or derived from N-(1-naphthalenyl)indole-3-carboxamide or
1072 1H-indol-(N-naphthyl)-3-carboxamide by substitution at the
1073 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
1074 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1075 cycloalkylmethyl, cycloalkylethyl,
1076 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1077 or 2-(4-morpholinyl)ethyl whether or not further substituted
1078 in the indole ring to any extent, whether or not substituted
1079 in the naphthyl ring to any extent.

1080 13. Any compound structurally analogous to, mimicking,
1081 or derived from N-(adamantan-1yl)indole-3-carboxamide or
1082 1H-indol-3-carboxamide-(1-adamantyl) by substitution at the
1083 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
1084 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1085 cycloalkylmethyl, cycloalkylethyl,
1086 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1087 or 2-(4-morpholinyl)ethyl whether or not further substituted
1088 in the indole ring to any extent.

1089 14. Any compound structurally analogous to, mimicking,
1090 or derived from N-(adamantan-1yl)indazole-3-carboxamide or
1091 1H-indazole-3-carboxamide-(1-adamantyl) by substitution at the
1092 nitrogen atom of the indazole ring by alkyl, alkyl halide,



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1093 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1094 cycloalkylmethyl, cycloalkylethyl,
1095 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1096 or 2-(4-morpholinyl)ethyl whether or not further substituted
1097 in the indazole ring to any extent.

1098 15. Any compound structurally analogous to, mimicking,
1099 or derived from
1100 N-[(1S)-1-(aminocarbonyl)-2-methylpropoyl]indazole-3-carboxyam
1101 ide or
1102 1H-indazole-3-carboxamide-N-[(1S)-1-(aminocarbonyl)-2-methylpr
1103 opoyl] by substitution at the nitrogen atom of the indazole
1104 ring by alkyl, alkyl halide, aryl halide, alkyl aryl halide,
1105 alkenyl, aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
1106 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1107 or 2-(4-morpholinyl)ethyl whether or not further substituted
1108 in the indazole ring to any extent.

1109 16. Any compound structurally analogous to, mimicking,
1110 or derived from 3-(1-naphthoyl)indazole or
1111 1H-indazole-3-yl-(1-naphthyl)methane by substitution at the
1112 nitrogen atom of the indazole ring by alkyl, alkyl halide,
1113 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
1114 cycloalkylmethyl, cycloalkylethyl,
1115 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1116 or 2-(4-morpholinyl)ethyl whether or not further substituted
1117 in the indazole ring to any extent, whether or not substituted
1118 in the naphthyl ring to any extent.

1119 17. Any compound structurally analogous to, mimicking,
1120 or derived from 3-(carboxylic acid 8-quinolinyl ester)indole



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1121 or 1H-indol-3-carboxylic acid-(8-quinolinyl)ester by
1122 substitution at the nitrogen atom of the indole ring by alkyl,
1123 alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
1124 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
1125 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
1126 or 2-(4-morpholinyl)ethyl whether or not further substituted
1127 in the indole ring to any extent, whether or not substituted
1128 in the quinoline ring to any extent.

1129 18. Any compound structurally related to
1130 2-(4-iodo-2,5-dimethoxyphenyl) ethanamine by substitution of
1131 the iodo moiety (4 position) with other halides, alkyl, alkyl
1132 halides, thioalkyl, cycloalkyl, cycloalkylhalides and/or
1133 substitution at the nitrogen atom of the ethanamine with
1134 alkyl, alkyl halide, alkenyl, cycloalkylmethyl,
1135 cycloalkylethyl, phenyl, benzyl whether or not further
1136 substituted in the (either) phenyl ring to any extent.

1137 19. Any compound structurally related to
1138 2,5-dimethoxy-4-chloroamphetamine by substitution of the
1139 chloro moiety (4 position) with other halides, alkyl, alkyl
1140 halides, thioalkyl, cycloalkyl, cycloalkylhalides and/or
1141 substitution at the nitrogen atom with alkyl, alkyl halide,
1142 alkenyl, cycloalkylmethyl, cycloalkylethyl, phenyl, benzyl
1143 whether or not further substituted in the (either) phenyl ring
1144 to any extent.

1145 20. Any compound structurally related to
1146 2-amino-1-phenyl-1-propanone (cathinone) by substitution of
1147 the amine with alkyl, alkyl halide, alkenyl, cycloalkylmethyl,
1148 cycloalkylethyl, phenyl, benzyl whether or not further



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1149 substituted in the (either) phenyl ring to any extent.

1150 21. Any compound structurally related to
1151 a-pyrrolidinopentiophenone (a-pvp) whether or not further
1152 substituted in the phenyl ring to any extent, whether or not
1153 further substituted in the pyrrolidine ring to any extent.

1154 b. A synthetic controlled substance or analogue in
1155 subdivision (4) or this subdivision does not include any of
1156 the following:

1157 1. Any substance for which there is an approved new
1158 drug application under the Federal Food, Drug, and Cosmetic
1159 Act.

1160 2. With respect to a particular person, any substance,
1161 if an exemption is in effect for investigational use, for that
1162 person, as provided by 21 U.S.C. § 355, and the person is
1163 registered as a controlled substance researcher as required
1164 under ~~section~~Section 152.12, ~~subdivision~~Subdivision 3, to the
1165 extent conduct with respect to the substance is pursuant to
1166 the exemption and registration.

1167 c. A controlled substance analogue is treated as a
1168 controlled substance in Schedule I.

1169 d. After the Alabama Department of Forensic Sciences
1170 has determined a substance to be a synthetic controlled
1171 substance analogue under this section, the department shall
1172 notify the Alabama Department of Public Health with
1173 information relevant to scheduling as provided by Section
1174 20-2-20."

1175 Section 2. Section 13A-12-214.4, Code of Alabama 1975,
1176 which prohibits the sale of psychoactive cannabinoids to



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1177 minors, is repealed.

1178 Section 3. This act shall become effective on October

1179 1, 2025.