

By: Representative Yancey

To: Drug Policy

HOUSE BILL NO. 1071

1 AN ACT TO AMEND SECTION 41-29-113, MISSISSIPPI CODE OF 1972,
 2 TO INCLUDE SIXTEEN SUBSTANCES AS SCHEDULE I CONTROLLED SUBSTANCES
 3 BECAUSE THESE DRUGS HAVE NO LEGITIMATE MEDICAL USE AND HAVE A HIGH
 4 POTENCY WITH GREAT POTENTIAL TO CAUSE HARM; TO AMEND SECTION
 5 41-29-119, MISSISSIPPI CODE OF 1972, TO INCLUDE
 6 SERDEXMETHYLPHENIDATE AS A SCHEDULE IV CONTROLLED SUBSTANCE
 7 BECAUSE THE DRUG HAS A CURRENTLY ACCEPTED MEDICAL USE AND A LOW
 8 POTENTIAL FOR ABUSE THAT MAY LEAD TO LIMITED PHYSICAL DEPENDENCE
 9 OR PSYCHOLOGICAL DEPENDENCE RELATIVE TO THE DRUGS OR OTHER
 10 SUBSTANCES IN SCHEDULE III; AND FOR RELATED PURPOSES.

11 BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF MISSISSIPPI:

12 **SECTION 1.** Section 41-29-113, Mississippi Code of 1972, is
 13 amended as follows:

14 41-29-113.

15 **SCHEDULE I**

16 (a) Schedule I consists of the drugs and other substances,
 17 by whatever official name, common or usual name, chemical name, or
 18 brand name designated, that is listed in this section.

19 (b) **Opiates.** Unless specifically excepted or unless listed
 20 in another schedule, any of the following opiates, including their
 21 isomers, esters, ethers, salts and salts of isomers, esters and



22 ethers, whenever the existence of these isomers, esters, ethers
23 and salts is possible within the specific chemical designation:

24 (1) Acetyl-alpha-methylfentanyl
25 (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);

26 (2) Acetylmethadol;

27 (3) Acetyl fentanyl
28 (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

29 (4) Acryl fentanyl
30 (N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide), also known as
31 acryloylfentanyl;

32 (5) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)
33 cyclohexylmethyl]benzamide);

34 (6) Allylprodine;

35 (7) Alphacetylmethadol, except levo-alphacetylmethadol
36 (levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM);

37 (8) Alphameprodine;

38 (9) Alphamethadol;

39 (10) Alpha-Methylfentanyl
40 (N-[1-alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide;
41 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine);

42 (11) Alpha-Methylthiofentanyl
43 (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide
44);

45 (12) Benzethidine;

46 (13) Betacetylmethadol;



47 (14) Beta-Hydroxyfentanyl
48 (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide);
49 (15) Beta-Hydroxy-3-methylfentanyl
50 (N-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidinyl]-N-phenylpr
51 opanamide);
52 (16) Beta-Hydroxythiofentanyl
53 (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpr
54 opionamide);
55 (17) Betameprodine;
56 (18) Betamethadol;
57 (19) Beta-Methyl fentanyl
58 (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide), also
59 known as β -methyl fentanyl;
60 (20) Beta'-Phenyl fentanyl
61 (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide), also
62 known as β' -phenyl fentanyl or 3-phenylpropanoyl fentanyl;
63 (21) Betaprodine;
64 (22) Butyrl fentanyl
65 (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide);
66 (23) Clonitazene;
67 (24) Crotonyl fentanyl
68 ((E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide);
69 (25) Cyclopentyl fentanyl
70 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);



71 (26) Cyclopropyl fentanyl
72 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);
73 (27) Dextromoramide;
74 (28) Diampromide;
75 (29) Diethylthiambutene;
76 (30) Difenoxylin;
77 (31) Dimenoxadol;
78 (32) Dimepheptanol;
79 (33) Dimethylthiambutene;
80 (34) Dioxaphetyl butyrate;
81 (35) Dipipanone;
82 (36) Ethylmethylthiambutene;
83 (37) Etonitazene;
84 (38) Etoxadoline;
85 (39) Fentanyl carbamate
86 (ethyl(1-phenethylpiperidin-4-yl)(phenyl)carbamate);
87 (40) Fentanyl-related substances, meaning any substance
88 not otherwise listed under another schedule and for which no
89 exemption or approval is in effect under Section 505 of the
90 Federal Food, Drug, and Cosmetic Act [21 USC 355] that is
91 structurally related to fentanyl by one or more of the following
92 modifications:
93 (A) Replacement of the phenyl portion of the
94 phenethyl group by any monocycle, whether or not further
95 substituted in or on the monocycle;



96 (B) Substitution in or on the phenethyl group with
97 alkyl, alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino or nitro
98 groups;

99 (C) Substitution in or on the piperidine ring with
100 alkyl, alkenyl, alkoxy, ester, ether, hydroxyl, halo, haloalkyl,
101 amino or nitro groups;

102 (D) Replacement of the aniline ring with any
103 aromatic monocycle whether or not further substituted in or on the
104 aromatic monocycle; and/or

105 (E) Replacement of the N-propionyl group by
106 another acyl group.

107 (41) 4-Fluoroisobutyryl fentanyl
108 (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide),
109 also known as para-fluoroisobutyryl fentanyl);

110 (42) 2'-Fluoro ortho-fluorofentanyl
111 (N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)
112 propionamide), also known as 2'-fluoro 2-fluorofentanyl;

113 (43) Furanyl fentanyl
114 (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide);

115 (44) Furethidine;

116 (45) Hydroxypethidine;

117 (46) Isobutyryl fentanyl
118 (N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide);

119 (47) Isotonitazene (N,N-diethyl-2-(2-(4
120 isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine);



121 (48) Isotonitazene
122 (N,N-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-y
123 l)ethan-1-amine);
124 (* * *49) Ketobemidone (including the optical and
125 geometric isomers);
126 (* * *50) Levomoramide;
127 (* * *51) Levophenacylmorphane;
128 (* * *52) Methoxyacetyl fentanyl
129 (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);
130 (* * *53) 4'-Methyl acetyl fentanyl
131 (N-(1-(4-methylphenethyl)piperidin-4-yl)-N-phenylacetamide);
132 (* * *54) 3-Methylfentanyl
133 (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);
134 (* * *55) 3-Methylthiofentanyl (N-[3-methyl-1-
135 (2-thienylethyl)-4-piperidinyl]-N-phenylpropanamide);
136 (* * *56) Morpheridine;
137 (* * *57) MPPP
138 (1-methyl-4-phenyl-4-propionoxypiperidine);
139 (* * *58) MT-45
140 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
141 (* * *59) Noracymethadol;
142 (* * *60) Norlevorphanol;
143 (* * *61) Normethadone;
144 (* * *62) Norpipanone;
145 (* * *63) Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-



146 (1-phenethylpiperidin-4-yl)acetamide);
147 (* * *64) Ortho-Fluoroacryl fentanyl
148 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide);
149 (* * *65) Ortho-Fluorobutyryl fentanyl
150 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide), also
151 known as 2-fluorobutyryl fentanyl;
152 (* * *66) Ortho-Fluorofentanyl
153 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide),
154 also known as 2-fluorofentanyl;
155 (* * *67) Ortho-Fluoroisobutyryl fentanyl
156 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);
157 (* * *68) Ortho-Methyl acetylfentanyl
158 (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide), also
159 known as 2-methyl acetylfentanyl;
160 (* * *69) Ortho-Methyl methoxyacetyl fentanyl
161 (2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)
162 acetamide), also known as 2-methyl methoxyacetyl fentanyl;
163 (* * *70) Para-Chloroisobutyryl fentanyl
164 (N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);
165 (* * *71) Para-Fluorobutyryl fentanyl
166 (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);
167 (* * *72) Para-Fluorofentanyl (N-(4-fluorophenyl)
168 -N-[1-(2-phenylethyl)-4-piperidinyl]propanamide);
169 (* * *73) Para-Fluoro furanyl fentanyl
170 N-(4-fluorophenyl)-N-



171 (1-phenethylpiperidin-4-yl) furan-2-carboxamide);
172 (* * *74) Para-Methoxybutyryl fentanyl
173 (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);
174 (* * *75) Para-Methylfentanyl
175 (N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide),
176 also known as 4-methylfentanyl);
177 (* * *76) PEPAP
178 (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
179 (* * *77) Phenadoxone;
180 (* * *78) Phenampromide;
181 (* * *79) Phenomorphan;
182 (* * *80) Phenoperidine;
183 (* * *81) Phenyl fentanyl
184 (N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide), also known as
185 benzoyl fentanyl;
186 (* * *82) Piritramide;
187 (* * *83) Proheptazine;
188 (* * *84) Properidine;
189 (* * *85) Propiram;
190 (* * *86) Racemoramide;
191 (* * *87) Tetrahydrofuranyl fentanyl
192 (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-
193 carboxamide);
194 (* * *88) Thiofentanyl
195 (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]propanamide);



196 (* * *89) Thiofuranyl fentanyl
197 (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide),
198 also known as 2-thiofuranyl fentanyl or thiophene fentanyl;

199 (* * *90) Tilidine;

200 (* * *91) Trimeperidine;

201 (* * *92) U-47700, (3,4-dichloro-N-
202 [2-(dimethylamino)cyclohexyl]-N-methylbenzamide);

203 (* * *93) Valeryl fentanyl
204 (N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide).

205 (c) **Opium derivatives.** Unless specifically excepted or
206 unless listed in another schedule, any of the following opium
207 derivatives, their salts, isomers and salts of isomers, whenever
208 the existence of these salts, isomers and salts of isomers is
209 possible within the specific chemical designation:

210 (1) Acetorphine;

211 (2) Acetyldihydrocodeine;

212 (3) Benzylmorphine;

213 (4) Codeine methylbromide;

214 (5) Codeine-N-Oxide;

215 (6) Cyrenorphine;

216 (7) Desomorphine;

217 (8) Dihydromorphine;

218 (9) Drotebanol;

219 (10) Etorphine (except hydrochloride salt);

220 (11) Heroin;



- 221 (12) Hydromorphinol;
- 222 (13) Methyldesorphine;
- 223 (14) Methyldihydromorphine;
- 224 (15) Monoacetylmorphine;
- 225 (16) Morphine methylbromide;
- 226 (17) Morphine methylsulfonate;
- 227 (18) Morphine-N-Oxide;
- 228 (19) Myrophine;
- 229 (20) Nicocodeine;
- 230 (21) Nicomorphine;
- 231 (22) Normorphine;
- 232 (23) Pholcodine;
- 233 (24) Thebacon.

234 (d) **Hallucinogenic substances.** Unless specifically excepted
235 or unless listed in another schedule, any material, compound,
236 mixture or preparation which contains any quantity of the
237 following substances, their salts, isomers (whether optical,
238 positional, or geometric) and salts of isomers, whenever the
239 existence of these salts, isomers and salts of isomers is possible
240 within the specific chemical designation:

- 241 (1) Alpha-ethyltryptamine;
- 242 (2) 4-bromo-2,5-dimethoxy-amphetamine;
- 243 (3) 4-bromo-2,5-dimethoxyphenethylamine;
- 244 (4) 2,5-dimethoxyamphetamine;
- 245 (5) 2,5-dimethoxy-4-ethylamphetamine (DOET);



- 246 (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine
247 (2C-T-7);
- 248 (7) 4-methoxyamphetamine;
- 249 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;
- 250 (9) 4-methyl-2,5-dimethoxy-amphetamine;
- 251 (10) 3,4-methylenedioxy amphetamine;
- 252 (11) 3,4-methylenedioxymethamphetamine (MDMA);
- 253 (12) 3,4-methylenedioxy-N-ethylamphetamine (also known
254 as N-ethyl-alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl
255 MDA, MDE, MDEA);
- 256 (13) N-hydroxy-3,4-methylenedioxyamphetamine (also
257 known as N-hydroxy MDA, N-OHMDA, and
258 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine);
- 259 (14) 3,4,5-trimethoxy amphetamine;
- 260 (15) 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
- 261 (16) Alpha-methyltryptamine (also known as AMT);
- 262 (17) Bufotenine;
- 263 (18) Diethyltryptamine;
- 264 (19) Dimethyltryptamine;
- 265 (20) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);
- 266 (21) Ibogaine;
- 267 (22) Lysergic acid diethylamide (LSD);
- 268 (23) (A) Marijuana (Hemp as defined and regulated
269 under Sections 69-25-201 through 69-25-221 and Cannabidiol
270 contained in a legend drug product approved by the Federal Food



271 and Drug Administration or obtained under Section 41-29-136 are
272 exempt under Schedule I);

273 (B) Hashish;

274 (24) Mescaline;

275 (25) Parahexyl;

276 (26) Peyote;

277 (27) N-ethyl-3-piperidyl benzilate;

278 (28) N-methyl-3-piperidyl benzilate;

279 (29) Psilocybin;

280 (30) Psilocyn;

281 (31) Tetrahydrocannabinols, meaning

282 tetrahydrocannabinols contained in a plant of the genus Cannabis
283 (cannabis plant), as well as the synthetic equivalents of the
284 substances contained in the cannabis plant, or in the resinous
285 extractives of such plant, and/or synthetic substances,
286 derivatives, and their isomers with similar chemical structure and
287 pharmacological activity to those substances contained in the
288 plant such as the following:

289 (A) 1 cis or trans tetrahydrocannabinol;

290 (B) 6 cis or trans tetrahydrocannabinol;

291 (C) 3,4 cis or trans tetrahydrocannabinol.

292 (Since nomenclature of these substances is not
293 internationally standardized, compounds of these structures,
294 regardless of atomic positions, are covered.)



295 ("Tetrahydrocannabinols" excludes dronabinol and nabilone.)
296 For purposes of this paragraph, tetrahydrocannabinols do not
297 include hemp or hemp products regulated under Sections 69-25-201
298 through 69-25-221.

299 However, the following products are exempted from control:

300 (i) THC-containing industrial products made
301 from cannabis stalks (e.g., paper, rope and clothing);

302 (ii) Processed cannabis plant materials used
303 for industrial purposes, such as fiber retted from cannabis stalks
304 for use in manufacturing textiles or rope;

305 (iii) Animal feed mixtures that contain
306 sterilized cannabis seeds and other ingredients (not derived from
307 the cannabis plant) in a formula designed, marketed and
308 distributed for nonhuman consumption;

309 (iv) Personal care products that contain oil
310 from sterilized cannabis seeds, such as shampoos, soaps, and body
311 lotions (if the products do not cause THC to enter the human
312 body);

313 (v) Hemp as regulated under Sections
314 69-25-201 through 69-25-221; and

315 (vi) Any product derived from the hemp plant
316 designed for human ingestion and/or consumption that is approved
317 by the United States Food and Drug Administration;

318 (32) Phencyclidine;

319 (33) Ethylamine analog of phencyclidine (PCE);



320 (34) Pyrrolidine analog of phencyclidine (PHP, PCPy);
321 (35) Thiophene analog of phencyclidine;
322 (36) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine (TCPy);
323 (37) 4-methylmethcathinone (mephedrone);
324 (38) 3,4-methylenedioxypropylvalerone (MDPV);
325 (39) 2-(2,5-dimethoxy-4-ethylphenyl)ethanamine (2C-E);
326 (40) 2-(2,5-dimethoxy-4-methylphenyl)ethanamine (2C-D);
327 (41) 2-(4-chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);
328 (42) 2-(4-iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);
329 or 2,5-dimethoxy-4-iodophenethylamine;
330 (43) 2-[4-(ethylthio)-2,5-dimethoxyphenyl]ethanamine
331 (2C-T-2);
332 (44)
333 2-[4-(isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);
334 (45) 2-(2,5-dimethoxyphenyl)ethanamine (2C-H);
335 (46) 2-(2,5-dimethoxy-4-nitro-phenyl)ethanamine (2C-N);
336 (47) 2-(2,5-dimethoxy-4-(n)-propylphenyl)ethanamine
337 (2C-P);
338 (48) 3,4-methylenedioxy-N-methylcathinone (methylone);
339 (49)
340 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine
341 (25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);
342 (50)
343 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine
344 (25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82);



345 (51)
346 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine or
347 N-[(2-methoxyphenyl)methyl]ethanamine (25I-NBOMe; 2C-I-NBOMe; 25I;
348 Cimbi-5);

349 (52) 7-bromo-5-(2-chlorophenyl)-1,3-dihydro-2H-1,
350 4-benzodiazepin-2-one (also known as Phenazepam);

351 (53) 7-(2-chlorophenyl)-4-ethyl-13-methyl-3-thia-1,8,
352 11,12-tetraazatricyclo[8.3.0.0]trideca-2(6),4,7,10,12-pentaene
353 (also known as Etizolam);

354 (54) *Salvia divinorum*;

355 (55) Synthetic cannabinoids. Unless specifically
356 excepted or unless listed in another schedule, any material,
357 compound, mixture, or preparation which contains any quantity of a
358 synthetic cannabinoid found in any of the following chemical
359 groups, whether or not substituted to any extent, or any of those
360 groups which contain any synthetic cannabinoid salts, isomers, or
361 salts of isomers, whenever the existence of such salts, isomers,
362 or salts of isomers is possible within the specific chemical
363 designation, including all synthetic cannabinoid chemical
364 analogues in such groups:

365 (A) (6aR,10aR)-9-(hydroxymethyl)-6,
366 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]
367 chromen-1-ol (also known as HU-210 or
368 1,1-dimethylheptyl-11-hydroxy-delta8-tetrahydrocannabinol);



369 (B) Naphthoylindoles and naphthylmethylinindoles,
370 being any compound structurally derived from 3-(1-naphthoyl)indole
371 or 1H-indol-3-yl-(1-naphthyl)methane, whether or not substituted
372 in the indole ring to any extent, or in the naphthyl ring to any
373 extent;

374 (C) Naphthoylpyrroles, being any compound
375 structurally derived from 3-(1-naphthoyl)pyrrole, whether or not
376 substituted in the pyrrole ring to any extent, or in the naphthyl
377 ring to any extent;

378 (D) Naphthylmethylinindenes, being any compound
379 structurally derived from 1-(1-naphthylmethyl)indene, whether or
380 not substituted in the indene ring to any extent or in the
381 naphthyl ring to any extent;

382 (E) Phenylacetylinindoles, being any compound
383 structurally derived from 3-phenylacetylinindole, whether or not
384 substituted in the indole ring to any extent or in the phenyl ring
385 to any extent;

386 (F) Cyclohexylphenols, being any compound
387 structurally derived from 2-(3-hydroxycyclohexyl)phenol, whether
388 or not substituted in the cyclohexyl ring to any extent or in the
389 phenolic ring to any extent;

390 (G) Benzoylinindoles, whether or not substituted in
391 the indole ring to any extent or in the phenyl ring to any extent;



392 (H) Adamantoylindoles, whether or not substituted
393 in the indole ring to any extent or in the adamantoyl ring system
394 to any extent;

395 (I) Tetrahydro derivatives of cannabinal and
396 3-alkyl homologues of cannabinal or of its tetrahydro derivatives,
397 except where contained in cannabis or cannabis resin;

398 (J) 3-Cyclopropylmethanone indole or
399 3-Cyclobutylmethanone indole or 3-Cyclopentylmethanone indole by
400 substitution at the nitrogen atom of the indole ring, whether or
401 not further substituted in the indole ring to any extent, whether
402 or not substituted on the cyclopropyl, cyclobutyl or cyclopentyl
403 rings to any extent;

404 (K) Quinoliny ester indoles, being any compound
405 structurally derived from 1H-indole-3carboxylic acid-8-quinoliny
406 ester, whether or not substituted in the indole ring to any extent
407 or the quinolone ring to any extent;

408 (L) 3-carboxamide-1H-indazoles, whether or not
409 substituted in the indazole ring to any extent and substituted to
410 any degree on the carboxamide nitrogen and
411 3-carboxamide-1H-indoles, whether or not substituted in the indole
412 ring to any extent and substituted to any degree on the
413 carboxamide nitrogen;

414 (M) Cycloalkanemethanone Indoles, whether or not
415 substituted at the nitrogen atom on the indole ring, whether or



416 not further substituted in the indole ring to any extent, whether
417 or not substituted on the cycloalkane ring to any extent;

418 (56) Naphthalen-1-yl

419 1-(5-fluoropentyl)-1H-indole-3-carboxylate, also known as NM2201
420 or CBL2201;

421 (57) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
422 pyrrolo[2,3-b]pyridine-3-carboxamide, also known as
423 5F-CUMYL-P7AICA or SGT-25;

424 (58) methyl

425 2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutano
426 ate, also known as 4F-MDMB-BINACA or 4F-MDMB-BUTINACA)

427 (* * * 59) 1-(4-methoxyphenyl)-N-methylpropan-2-amine,
428 also known as para-methoxymethamphetamine or PMMA

429 (60) ethyl 2-(1-(5-fluoropentyl)

430 -1H-indazole-3-carboxamido)-3,3-dimethylbutanoate, also known as
431 5F-EDMB-PINACA;

432 (61) methyl

433 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,

434 3-dimethylbutanoate, also known as 5F-MDMB-PICA or 5F-MDMB-2201;

435 (62)

436 N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide,

437 also known as FUB-AKB48 or FUB-APINACA or AKB48

438 N-(4-fluorobenzyl);



439 (63)
440 (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)
441 methanone, also known as FUB-144;

442 (64) N-ethylhexedrone, also known as
443 α -ethylaminohexanophenone or 2-(ethylamino)-1-phenylhexan-1-one;

444 (65) alpha-pyrrolidinohexanophenone, also known as
445 α -PHP or α -pyrrolidinohexanophenone or
446 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one);

447 (66) 4-methyl-alpha-ethylaminopentiophenone, also known
448 as 4-MEAP or 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one);

449 (67) 4'-methyl-alpha-pyrrolidinohexiophenone, also
450 known as MPHP or 4'-methyl-alpha-pyrrolidinohexanophenone or
451 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);

452 (68) alpha-pyrrolidinoheptaphenone (also known as PV8;
453 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one);

454 (69) 4'-chloro-alpha-pyrrolidinovalerophenone, also
455 known as 4-chloro- α -PVP or 4'-chloro- α -pyrrolidinopentiophenone or
456 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one);

457 (70)
458 2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one, also known as
459 methoxetamine or MXE.

460 (e) **Depressants.** Unless specifically excepted or unless
461 listed in another schedule, any material, compound, mixture, or
462 preparation which contains any quantity of the following
463 substances having a depressant effect on the central nervous



464 system, including their salts, isomers, and salts of isomers,
465 whenever the existence of such salts, isomers, and salts of
466 isomers is possible within the specific chemical designation:

467 (1) Clonazolam,
468 6-(2-chlorophenyl)-1-methyl-8-nitro-4H-[1,2,4]triazolo[4,3-a][1,4]
469 benzodiazepine;

470 (2) Flualprazolam,
471 8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4
472]benzodiazepine;

473 (3) Flubromazepam,
474 7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one;

475 (4) Flubromazolam,
476 8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]
477 benzodiazepine;

478 (5) Gamma-hydroxybutyric acid (other names include:
479 GHB, gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic
480 acid; sodium oxybate; sodium oxybutyrate);

481 (6) Mecloqualone;

482 (7) Methaqualone.

483 (f) **Stimulants.** Any material, compound, mixture or
484 preparation which contains any quantity of the following central
485 nervous system stimulants including optical salts, isomers and
486 salts of isomers unless specifically excepted or unless listed in
487 another schedule:

488 (1) Aminorex;



489 (2) N-benzylpiperazine (also known as BZP and
490 1-benzylpiperazine);
491 (3) Cathinone;
492 (4) 4,4'-Dimethylaminorex, also known as 4,4'-DMAR or
493 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine;
494 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine);
495 (* * *5) Fenethylamine;
496 (* * *6) Methcathinone;
497 (* * *7) 4-methylaminorex (also known as
498 2-amino-4-methyl-5-phenyl-2-oxazoline);
499 (* * *8) N-ethylamphetamine;
500 (* * *9) Any material, compound, mixture or
501 preparation which contains any quantity of
502 N,N-dimethylamphetamine. (Other names include:
503 N,N,-alpha-trimethyl-benzeneethanamine and
504 N,N-alpha-trimethylphenethylamine);
505 (* * *10) Synthetic cathinones. (A) Unless listed in
506 another schedule, any compound other than bupropion that is
507 structurally derived from 2-Amino-1-phenyl-1-propanone by
508 modification in any of the following ways:
509 (i) By substitution in the phenyl ring to any
510 extent with alkyl, alkoxy, alkylendioxy, haloalkyl or halide
511 substituents, whether or not further substituted in the phenyl
512 ring by one or more other univalent substituents;



513 (ii) By substitution at the 3-position with
514 an alkyl substituent;

515 (iii) By substitution at the nitrogen atom
516 with alkyl or dialkyl groups, or by inclusion of the nitrogen atom
517 in a cyclic structure.

518 (B) The compounds covered in this paragraph
519 (* * * 10) include, but are not limited to, any material,
520 compound, mixture or preparation which contains any quantity of a
521 synthetic cathinone found in any of the following compounds,
522 whether or not substituted to any extent, or any of these
523 compounds which contain any synthetic cathinone, or salts,
524 isomers, or salts of isomers, whenever the existence of such
525 salts, isomers or salts of isomers is possible, unless
526 specifically excepted or listed in another schedule:

527 (i) 4-methyl-N-ethylcathinone ("4-MEC");

528 (ii) 4-methyl-alpha-pyrrolidinopropiophenone
529 ("4-MePPP");

530 (iii) Alpha-pyrrolidinopentiophenone
531 ("α-PVP");

532 (iv) 1-(1,3-benzodioxol-5-yl)-2-
533 (methylamino)butan-1-one ("butylone");

534 (v) 2-(methylamino)-1-phenylpentan-1-one
535 ("pentedrone");

536 (vi) 1-(1,3-benzodioxol-5-yl)-2-
537 (methylamino)pentan-1-one ("pentylone");



538 (vii) 4-fluoro-N-methylcathinone ("4-FMC");
539 (viii) 3-fluoro-N-methylcathinone ("3-FMC");
540 (ix) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)
541 pentan-1-one ("naphyrone");
542 (x) Alpha-pyrrolidinobutiophenone
543 ("α-PBP"); * * *
544 (xi) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)
545 -pentan-1-one (N-ethylpentylone, ephylone) * * *; and
546 (xii)
547 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one, also know as
548 N-ethylpentylone.

549 **SECTION 2.** Section 41-29-119, Mississippi Code of 1972, is
550 amended as follows:

551 41-29-119. (A) The controlled substances listed in this
552 section are included in Schedule IV.

553 **SCHEDULE IV**

554 (a) **Narcotic drugs.** Unless specifically excepted or
555 unless listed in another schedule, any material, compound, mixture
556 or preparation which contains limited quantities of the following
557 narcotic drugs, or any salts thereof:

558 (1) Not more than one (1) milligram of difenoxin
559 and not less than twenty-five (25) micrograms of atropine sulfate
560 per dosage unit;



561 (2) Dextropropoxyphene, including its salts
562 (Darvon, Darvon-N; also found in Darvon compound and Darvocet-N,
563 etc.);

564 (3)
565 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its
566 salts, optical and geometric isomers and salts of these isomers
567 (including tramadol).

568 (b) **Depressants.** Any material, compound, mixture or
569 preparation which contains any quantity of the following
570 substances:

571 (1) Alfaxalone;

572 (2) Alprazolam;

573 (3) Barbital;

574 (4) Brexanolone;

575 (5) Bromazepam;

576 (6) Camazepam;

577 (7) Carisoprodol;

578 (8) Chloral betaine;

579 (9) Chloral hydrate;

580 (10) Chlordiazepoxide and its salts, but does not
581 include chlordiazepoxide hydrochloride and clidinium bromide or
582 chlordiazepoxide and esterified estrogens;

583 (11) Clobazam;

584 (12) Clonazepam;

585 (13) Clorazepate;



- 586 (14) Clotiazepam;
587 (15) Cloxazolam;
588 (16) Delorazepam;
589 (17) Diazepam;
590 (18) Dichloralphenazone;
591 (19) Estazolam;
592 (20) Ethchlorvynol;
593 (21) Ethinamate;
594 (22) Ethyl loflazepate;
595 (23) Fludiazepam;
596 (24) Flunitrazepam;
597 (25) Flurazepam;
598 (26) Fospropofol;
599 (27) Halazepam;
600 (28) Haloxazolam;
601 (29) Ketazolam;
602 (30) Lemborexant;
603 (31) Loprazolam;
604 (32) Lorazepam;
605 (33) Lormetazepam;
606 (34) Mebutamate;
607 (35) Medazepam;
608 (36) Meprobamate;
609 (37) Methohexital;
610 (38) Methylphenobarbital;



- 611 (39) Midazolam;
612 (40) Nimetazepam;
613 (41) Nitrazepam;
614 (42) Nordiazepam;
615 (43) Oxazepam;
616 (44) Oxazolam;
617 (45) Paraldehyde;
618 (46) Petrichloral;
619 (47) Phenobarbital;
620 (48) Pinazepam;
621 (49) Prazepam;
622 (50) Quazepam;
623 (51) Remimazolam;
624 (52) Suvorexant;
625 (53) Temazepam;
626 (54) Tetrazepam;
627 (55) Triazolam;
628 (56) Zaleplon;
629 (57) Zolpidem;
630 (58) Zopiclone.

631 (c) **Fenfluramine.**

632 (d) **Lorcaserin.** Any material, compound, mixture, or
633 preparation which contains any quantity of Lorcaserin, including
634 its salts, isomers, and salts of such isomers, whenever the



635 existence of such salts, isomers, and salts of isomers is
636 possible.

637 (e) **Stimulants.** Any material, compound, mixture or
638 preparation which contains any quantity of the following
639 substances:

- 640 (1) Cathine ((+/-) Norpseudoephedrine);
641 (2) Diethylpropion;
642 (3) Fencamfamin;
643 (4) Fenproporex;
644 (5) Mazindol;
645 (6) Mefenorex;
646 (7) Modafinil;
647 (8) Pemoline (including any organometallic
648 complexes and chelates thereof);
649 (9) Phentermine;
650 (10) Pipradrol;
651 (11) Serdexmethylphenidate;
652 (* * *12) Sibutramine;
653 (* * *13) Solriamfetol;
654 (* * *14) SPA
655 ((-)-1-dimethylamino-1,2-diphenylethane).

656 (f) **Other substances.**

- 657 (1) Pentazocine;
658 (2) Butorphanol (including its optical isomers);



659 (3) Eluxadoline
660 (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopr
661 opyl][(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-meth
662 oxybenzoic acid); (including its optical isomers) and its salts,
663 isomers, and salts of isomers.

664 (B) Any material, compound, mixture or preparation which
665 contains any quantity of a Schedule IV controlled substance and is
666 listed as an exempt substance in 21 CFR, Section 1308.22, 1308.24,
667 1308.26, 1308.32 or 1308.34, shall be exempted from the provisions
668 of the Uniform Controlled Substances Law.

669 **SECTION 3.** This act shall take effect and be in force from
670 and after July 1, 2023.

