

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  $\beta$ -methyl fentanyl;  
60 (20) Beta'-Phenyl fentanyl  
61 (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide), also  
62 known as  $\beta'$ -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) Etoxadine;  
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, hydroxy, halo, haloalkyl, amino or nitro  
98 groups;

99 (C) Substitution in or on the piperidine ring with  
100 alkyl, alkenyl, alkoxy, ester, ether, hydroxy, 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    $\alpha$ -ethylaminohexanophenone or 2-(ethylamino)-1-phenylhexan-1-one;

444                   (65) alpha-pyrrolidinohexanophenone, also known as  
445    $\alpha$ -PHP or  $\alpha$ -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- $\alpha$ -PVP or 4'-chloro- $\alpha$ -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 benzodiazepin;

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.

