

By: Representative Yancey

To: Drug Policy

HOUSE BILL NO. 232

1 AN ACT TO AMEND SECTION 41-29-113, MISSISSIPPI CODE OF 1972,  
 2 TO INCLUDE 20 FENTANYL-RELATED SUBSTANCES, MT-45, NM2201,  
 3 5F-CUMYL-P7AICA AND PMMA AS SCHEDULE I CONTROLLED SUBSTANCES  
 4 BECAUSE THESE DRUGS HAVE NO LEGITIMATE MEDICAL USE AND HAVE A HIGH  
 5 POTENCY WITH GREAT POTENTIAL TO CAUSE HARM; TO ALPHABETIZE CERTAIN  
 6 SUBSTANCES TO CONFORM THE LIST OF SCHEDULE I OPIATES TO THE CODE  
 7 OF FEDERAL REGULATIONS; TO AMEND SECTION 41-29-115, MISSISSIPPI  
 8 CODE OF 1972, TO INCLUDE OLICERIDINE AS A SCHEDULE II CONTROLLED  
 9 SUBSTANCE BECAUSE OLICERIDINE HAS A CURRENTLY ACCEPTED MEDICAL USE  
 10 BUT HAS A HIGH POTENTIAL FOR ABUSE THAT MAY LEAD TO SEVERE  
 11 PSYCHOLOGICAL OR PHYSICAL DEPENDENCE; TO AMEND SECTION 41-29-119,  
 12 MISSISSIPPI CODE OF 1972, TO INCLUDE LEMBOREXANT AND REMIMAZOLAM  
 13 AS SCHEDULE IV CONTROLLED SUBSTANCES BECAUSE THESE DRUGS HAVE A  
 14 CURRENTLY ACCEPTED MEDICAL USE AND A LOW POTENTIAL FOR ABUSE THAT  
 15 MAY LEAD TO LIMITED PHYSICAL DEPENDENCE OR PSYCHOLOGICAL  
 16 DEPENDENCE RELATIVE TO THE DRUGS OR OTHER SUBSTANCES IN SCHEDULE  
 17 III; TO AMEND SECTION 41-29-121, MISSISSIPPI CODE OF 1972, TO  
 18 INCLUDE CENOBAMATE AND LASMIDITAN AS SCHEDULE V CONTROLLED  
 19 SUBSTANCES BECAUSE THESE DRUGS HAVE A CURRENTLY ACCEPTED MEDICAL  
 20 USE AND A LOW POTENTIAL FOR ABUSE THAT MAY LEAD TO LIMITED  
 21 PHYSICAL DEPENDENCE OR PSYCHOLOGICAL DEPENDENCE RELATIVE TO THE  
 22 DRUGS OR OTHER SUBSTANCES IN SCHEDULE IV; AND FOR RELATED  
 23 PURPOSES.

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

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

27 41-29-113.

28 **SCHEDULE I**



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

32 (b) **Opiates.** Unless specifically excepted or unless listed  
33 in another schedule, any of the following opiates, including their  
34 isomers, esters, ethers, salts and salts of isomers, esters and  
35 ethers, whenever the existence of these isomers, esters, ethers  
36 and salts is possible within the specific chemical designation:

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

39 (2) Acetylmethadol;

40 ( \* \* \*3) Acetyl fentanyl  
41 (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

42 (4) Acryl fentanyl  
43 (N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide), also known as  
44 acryloylfentanyl;

45 ( \* \* \*5) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)  
46 cyclohexylmethyl]benzamide);

47 \* \* \*

48 ( \* \* \*6) Allylprodine;

49 ( \* \* \*7) Alphacetylmethadol, except  
50 levo-alphacetylmethadol (levo-alpha-acetylmethadol, levomethadyl  
51 acetate, or LAAM);

52 ( \* \* \*8) Alphameprodine;

53 ( \* \* \*9) Alphamethadol;



54 ( \* \* \*10) Alpha-Methylfentanyl  
55 (N-[1-alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide;  
56 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine);  
57 ( \* \* \*11) Alpha-Methylthiofentanyl  
58 (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide  
59 );  
60 ( \* \* \*12) Benzethidine;  
61 ( \* \* \*13) Betacetylmethadol;  
62 ( \* \* \*14) Beta-Hydroxyfentanyl  
63 (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide);  
64 ( \* \* \*15) Beta-Hydroxy-3-methylfentanyl  
65 (N-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidinyl]-N-phenylpr  
66 opamide);  
67 (16) Beta-Hydroxythiofentanyl  
68 (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpr  
69 opionamide);  
70 ( \* \* \*17) Betameprodine;  
71 ( \* \* \*18) Betamethadol;  
72 (19) Beta-Methyl fentanyl  
73 (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide), also  
74 known as  $\beta$ -methyl fentanyl;  
75 (20) Beta'-Phenyl fentanyl  
76 (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide), also  
77 known as  $\beta'$ -phenyl fentanyl or 3-phenylpropanoyl fentanyl;  
78 ( \* \* \*21) Betaprodine;



79 ( \* \* \*22) Butyrl fentanyl  
80 (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide);  
81 ( \* \* \*23) Clonitazene;  
82 (24) Crotonyl fentanyl  
83 ((E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide);  
84 (25) Cyclopentyl fentanyl  
85 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);  
86 (26) Cyclopropyl fentanyl  
87 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);  
88 ( \* \* \*27) Dextromoramide;  
89 ( \* \* \*28) Diampromide;  
90 ( \* \* \*29) Diethylthiambutene;  
91 ( \* \* \*30) Difenoquin;  
92 ( \* \* \*31) Dimenoxadol;  
93 ( \* \* \*32) Dimepheptanol;  
94 ( \* \* \*33) Dimethylthiambutene;  
95 ( \* \* \*34) Dioxaphetyl butyrate;  
96 ( \* \* \*35) Dipipanone;  
97 ( \* \* \*36) Ethylmethylthiambutene;  
98 ( \* \* \*37) Etonitazene;  
99 ( \* \* \*38) Etoxadine;  
100 (39) Fentanyl carbamate  
101 (ethyl(1-phenethylpiperidin-4-yl)(phenyl)carbamate);  
102 ( \* \* \*40) Fentanyl-related substances, meaning any  
103 substance not otherwise listed under another schedule and for



104 which no exemption or approval is in effect under Section 505 of  
105 the Federal Food, Drug, and Cosmetic Act [21 USC 355] that is  
106 structurally related to fentanyl by one or more of the following  
107 modifications:

108 (A) Replacement of the phenyl portion of the  
109 phenethyl group by any monocycle, whether or not further  
110 substituted in or on the monocycle;

111 (B) Substitution in or on the phenethyl group with  
112 alkyl, alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino or nitro  
113 groups;

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

117 (D) Replacement of the aniline ring with any  
118 aromatic monocycle whether or not further substituted in or on the  
119 aromatic monocycle; and/or

120 (E) Replacement of the N-propionyl group by  
121 another acyl group.

122 \* \* \*

123 (41) 4-Fluoroisobutyryl fentanyl  
124 (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide),  
125 also known as para-fluoroisobutyryl fentanyl);

126 (42) 2'-Fluoro ortho-fluorofentanyl  
127 (N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)  
128 propionamide), also known as 2'-fluoro 2-fluorofentanyl;



129 (43) Furanyl fentanyl  
130 (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide);  
131 ( \* \* \*44) Furethidine;  
132 ( \* \* \*45) Hydroxypethidine;  
133 (46) Isobutyryl fentanyl  
134 (N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide);  
135 ( \* \* \*47) Isotonitazene (N,N-diethyl-2-(2-(4  
136 isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine);  
137 ( \* \* \*48) Ketobemidone (including the optical and  
138 geometric isomers);  
139 ( \* \* \*49) Levomoramide;  
140 ( \* \* \*50) Levophenacetylmorphan;  
141 (51) Methoxyacetyl fentanyl  
142 (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);  
143 (52) 4'-Methyl acetyl fentanyl  
144 (N-(1-(4-methylphenethyl)piperidin-4-yl)-N-phenylacetamide);  
145 ( \* \* \*53) 3-Methylfentanyl  
146 (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);  
147 ( \* \* \*54) 3-Methylthiofentanyl (N-[3-methyl-1-  
148 (2-thienylethyl)-4-piperidinyl]-N-phenylpropanamide);  
149 ( \* \* \*55) Morpheridine;  
150 ( \* \* \*56) MPPP  
151 (1-methyl-4-phenyl-4-propionoxypiperidine);  
152 ( \* \* \*57) \* \* \* MT-45  
153 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);



154 ( \* \* \*58) Noracymethadol;  
155 ( \* \* \*59) Norlevorphanol;  
156 ( \* \* \*60) Normethadone;  
157 ( \* \* \*61) Norpipanone;  
158 (62) Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-  
159 (1-phenethylpiperidin-4-yl)acetamide);  
160 (63) Ortho-Fluoroacryl fentanyl  
161 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide);  
162 (64) Ortho-Fluorobutyryl fentanyl  
163 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide), also  
164 known as 2-fluorobutyryl fentanyl;  
165 (65) Ortho-Fluorofentanyl  
166 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide),  
167 also known as 2-fluorofentanyl;  
168 (66) Ortho-Fluoroisobutyryl fentanyl  
169 (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);  
170 (67) Ortho-Methyl acetylfentanyl  
171 (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide), also  
172 known as 2-methyl acetylfentanyl;  
173 (68) Ortho-Methyl methoxyacetyl fentanyl  
174 (2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)  
175 acetamide), also known as 2-methyl methoxyacetyl fentanyl;  
176 (69) Para-Chloroisobutyryl fentanyl  
177 (N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);



178                   (70) Para-Fluorobutyryl fentanyl  
179 (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);  
180                   ( \* \* \*71)   Para-Fluorofentanyl (N-(4-fluorophenyl)  
181 -N-[1-(2-phenylethyl)-4-piperidinyl]propanamide);  
182                   (72) Para-Fluoro furanyl fentanyl N-(4-fluorophenyl)-N-  
183 (1-phenethylpiperidin-4-yl)furan-2-carboxamide);  
184                   (73) Para-Methoxybutyryl fentanyl  
185 (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);  
186                   (74) Para-Methylfentanyl  
187 (N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide),  
188 also known as 4-methylfentanyl);  
189                   ( \* \* \*75)   PEPAP  
190 (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);  
191                   ( \* \* \*76)   Phenadoxone;  
192                   ( \* \* \*77)   Phenampromide;  
193                   ( \* \* \*78)   Phenomorphane;  
194                   ( \* \* \*79)   Phenoperidine;  
195                   (80) Phenyl fentanyl  
196 (N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide), also known as  
197 benzoyl fentanyl;  
198                   ( \* \* \*81)   Piritramide;  
199                   ( \* \* \*82)   Proheptazine;  
200                   ( \* \* \*83)   Propoperidine;  
201                   ( \* \* \*84)   Propiram;  
202                   ( \* \* \*85)   Racemoramide;





203                   (86) Tetrahydrofuranyl fentanyl  
204 (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-  
205 carboxamide);  
206                   ( \* \* \*87) Thiofentanyl  
207 (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]propanamide);  
208                   (88) Thiofuranyl fentanyl  
209 (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide),  
210 also known as 2-thiofuranyl fentanyl or thiophene fentanyl;  
211                   ( \* \* \*89) Tilidine;  
212                   ( \* \* \*90) Trimeperidine;  
213                   ( \* \* \*91) U-47700, (3,4-dichloro-N-  
214 [2-(dimethylamino)cyclohexyl]-N-methylbenzamide) \* \* \*;  
215                   (92) Valeryl fentanyl  
216 (N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide).

217           (c) **Opium derivatives.** Unless specifically excepted or  
218 unless listed in another schedule, any of the following opium  
219 derivatives, their salts, isomers and salts of isomers, whenever  
220 the existence of these salts, isomers and salts of isomers is  
221 possible within the specific chemical designation:

- 222                   (1) Acetorphine;
- 223                   (2) Acetyldihydrocodeine;
- 224                   (3) Benzylmorphine;
- 225                   (4) Codeine methylbromide;
- 226                   (5) Codeine-N-Oxide;
- 227                   (6) Cyprenorphine;



- 228 (7) Desomorphine;
- 229 (8) Dihydromorphine;
- 230 (9) Drotebanol;
- 231 (10) Etorphine (except hydrochloride salt);
- 232 (11) Heroin;
- 233 (12) Hydromorphenol;
- 234 (13) Methyldesorphine;
- 235 (14) Methyldihydromorphine;
- 236 (15) Monoacetylmorphine;
- 237 (16) Morphine methylbromide;
- 238 (17) Morphine methylsulfonate;
- 239 (18) Morphine-N-Oxide;
- 240 (19) Myrophine;
- 241 (20) Nicocodeine;
- 242 (21) Nicomorphine;
- 243 (22) Normorphine;
- 244 (23) Pholcodine;
- 245 (24) Thebacon.

246 (d) **Hallucinogenic substances.** Unless specifically excepted  
247 or unless listed in another schedule, any material, compound,  
248 mixture or preparation which contains any quantity of the  
249 following substances, their salts, isomers (whether optical,  
250 positional, or geometric) and salts of isomers, whenever the  
251 existence of these salts, isomers and salts of isomers is possible  
252 within the specific chemical designation:



- 253 (1) Alpha-ethyltryptamine;
- 254 (2) 4-bromo-2,5-dimethoxy-amphetamine;
- 255 (3) 4-bromo-2,5-dimethoxyphenethylamine;
- 256 (4) 2,5-dimethoxyamphetamine;
- 257 (5) 2,5-dimethoxy-4-ethylamphetamine (DOET);
- 258 (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine
- 259 (2C-T-7);
- 260 (7) 4-methoxyamphetamine;
- 261 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;
- 262 (9) 4-methyl-2,5-dimethoxy-amphetamine;
- 263 (10) 3,4-methylenedioxy amphetamine;
- 264 (11) 3,4-methylenedioxymethamphetamine (MDMA);
- 265 (12) 3,4-methylenedioxy-N-ethylamphetamine (also known
- 266 as N-ethyl-alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl
- 267 MDA, MDE, MDEA);
- 268 (13) N-hydroxy-3,4-methylenedioxyamphetamine (also
- 269 known as N-hydroxy MDA, N-OHMDA, and
- 270 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine);
- 271 (14) 3,4,5-trimethoxy amphetamine;
- 272 (15) 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
- 273 (16) Alpha-methyltryptamine (also known as AMT);
- 274 (17) Bufotenine;
- 275 (18) Diethyltryptamine;
- 276 (19) Dimethyltryptamine;
- 277 (20) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);



- 278 (21) Ibogaine;
- 279 (22) Lysergic acid diethylamide (LSD);
- 280 (23) (A) Marijuana (Hemp as defined and regulated
- 281 under Sections 69-25-201 through 69-25-221 and Cannabidiol
- 282 contained in a legend drug product approved by the Federal Food
- 283 and Drug Administration or obtained under Section 41-29-136 are
- 284 exempt under Schedule I);
- 285 (B) Hashish;
- 286 (24) Mescaline;
- 287 (25) Parahexyl;
- 288 (26) Peyote;
- 289 (27) N-ethyl-3-piperidyl benzilate;
- 290 (28) N-methyl-3-piperidyl benzilate;
- 291 (29) Psilocybin;
- 292 (30) Psilocyn;
- 293 (31) Tetrahydrocannabinols, meaning
- 294 tetrahydrocannabinols contained in a plant of the genus Cannabis
- 295 (cannabis plant), as well as the synthetic equivalents of the
- 296 substances contained in the cannabis plant, or in the resinous
- 297 extractives of such plant, and/or synthetic substances,
- 298 derivatives, and their isomers with similar chemical structure and
- 299 pharmacological activity to those substances contained in the
- 300 plant such as the following:
- 301 (A) 1 cis or trans tetrahydrocannabinol;
- 302 (B) 6 cis or trans tetrahydrocannabinol;



303 (C) 3,4 cis or trans tetrahydrocannabinol.  
304 (Since nomenclature of these substances is not  
305 internationally standardized, compounds of these structures,  
306 regardless of atomic positions, are covered.)  
307 ("Tetrahydrocannabinols" excludes dronabinol and nabilone.)  
308 For purposes of this paragraph, tetrahydrocannabinols do not  
309 include hemp or hemp products regulated under Sections 69-25-201  
310 through 69-25-221.

311 However, the following products are exempted from control:

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

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

317 (iii) Animal feed mixtures that contain  
318 sterilized cannabis seeds and other ingredients (not derived from  
319 the cannabis plant) in a formula designed, marketed and  
320 distributed for nonhuman consumption;

321 (iv) Personal care products that contain oil  
322 from sterilized cannabis seeds, such as shampoos, soaps, and body  
323 lotions (if the products do not cause THC to enter the human  
324 body);

325 (v) Hemp as regulated under Sections  
326 69-25-201 through 69-25-221; and



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

330 (32) Phencyclidine;

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

332 (34) Pyrrolidine analog of phencyclidine (PHP, PCPy);

333 (35) Thiophene analog of phencyclidine;

334 (36) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine (TCPy);

335 (37) 4-methylmethcathinone (mephedrone);

336 (38) 3,4-methylenedioxypropylamphetamine (MDPV);

337 (39) 2-(2,5-dimethoxy-4-ethylphenyl)ethanamine (2C-E);

338 (40) 2-(2,5-dimethoxy-4-methylphenyl)ethanamine (2C-D);

339 (41) 2-(4-chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

340 (42) 2-(4-iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

341 or 2,5-dimethoxy-4-iodophenethylamine;

342 (43) 2-[4-(ethylthio)-2,5-dimethoxyphenyl]ethanamine

343 (2C-T-2);

344 (44)

345 2-[4-(isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

346 (45) 2-(2,5-dimethoxyphenyl)ethanamine (2C-H);

347 (46) 2-(2,5-dimethoxy-4-nitro-phenyl)ethanamine (2C-N);

348 (47) 2-(2,5-dimethoxy-4-(n)-propylphenyl)ethanamine

349 (2C-P);

350 (48) 3,4-methylenedioxy-N-methylcathinone (methylone);



351 (49)  
352 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine  
353 (25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);  
354 (50)  
355 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine  
356 (25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82);  
357 (51)  
358 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine or  
359 N-[(2-methoxyphenyl)methyl]ethanamine (25I-NBOMe; 2C-I-NBOMe; 25I;  
360 Cimbi-5);  
361 (52) 7-bromo-5-(2-chlorophenyl)-1,3-dihydro-2H-1,  
362 4-benzodiazepin-2-one (also known as Phenazepam);  
363 (53) 7-(2-chlorophenyl)-4-ethyl-13-methyl-3-thia-1,8,  
364 11,12-tetraazatricyclo[8.3.0.0]trideca-2(6),4,7,10,12-pentaene  
365 (also known as Etizolam);  
366 (54) Salvia divinorum;  
367 (55) Synthetic cannabinoids. Unless specifically  
368 excepted or unless listed in another schedule, any material,  
369 compound, mixture, or preparation which contains any quantity of a  
370 synthetic cannabinoid found in any of the following chemical  
371 groups, whether or not substituted to any extent, or any of those  
372 groups which contain any synthetic cannabinoid salts, isomers, or  
373 salts of isomers, whenever the existence of such salts, isomers,  
374 or salts of isomers is possible within the specific chemical



375 designation, including all synthetic cannabinoid chemical  
376 analogues in such groups:

377 (A) (6aR,10aR)-9-(hydroxymethyl)-6,  
378 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]  
379 chromen-1-ol (also known as HU-210 or  
380 1,1-dimethylheptyl-11-hydroxy-delta8-tetrahydrocannabinol);

381 (B) Naphthoylindoles and naphthylmethylindoles,  
382 being any compound structurally derived from 3-(1-naphthoyl)indole  
383 or 1H-indol-3-yl-(1-naphthyl)methane, whether or not substituted  
384 in the indole ring to any extent, or in the naphthyl ring to any  
385 extent;

386 (C) Naphthoylpyrroles, being any compound  
387 structurally derived from 3-(1-naphthoyl)pyrrole, whether or not  
388 substituted in the pyrrole ring to any extent, or in the naphthyl  
389 ring to any extent;

390 (D) Naphthylmethylindenes, being any compound  
391 structurally derived from 1-(1-naphthylmethyl)indene, whether or  
392 not substituted in the indene ring to any extent or in the  
393 naphthyl ring to any extent;

394 (E) Phenylacetylindoles, being any compound  
395 structurally derived from 3-phenylacetylindole, whether or not  
396 substituted in the indole ring to any extent or in the phenyl ring  
397 to any extent;

398 (F) Cyclohexylphenols, being any compound  
399 structurally derived from 2-(3-hydroxycyclohexyl)phenol, whether





400 or not substituted in the cyclohexyl ring to any extent or in the  
401 phenolic ring to any extent;

402 (G) Benzoylindoles, whether or not substituted in  
403 the indole ring to any extent or in the phenyl ring to any extent;

404 (H) Adamantoylindoles, whether or not substituted  
405 in the indole ring to any extent or in the adamantoyl ring system  
406 to any extent;

407 (I) Tetrahydro derivatives of cannabinal and  
408 3-alkyl homologues of cannabinal or of its tetrahydro derivatives,  
409 except where contained in cannabis or cannabis resin;

410 (J) 3-Cyclopropylmethanone indole or  
411 3-Cyclobutylmethanone indole or 3-Cyclopentylmethanone indole by  
412 substitution at the nitrogen atom of the indole ring, whether or  
413 not further substituted in the indole ring to any extent, whether  
414 or not substituted on the cyclopropyl, cyclobutyl or cyclopentyl  
415 rings to any extent;

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

420 (L) 3-carboxamide-1H-indazoles, whether or not  
421 substituted in the indazole ring to any extent and substituted to  
422 any degree on the carboxamide nitrogen and  
423 3-carboxamide-1H-indoles, whether or not substituted in the indole



424 ring to any extent and substituted to any degree on the  
425 carboxamide nitrogen;

426 (M) Cycloalkanemethanone Indoles, whether or not  
427 substituted at the nitrogen atom on the indole ring, whether or  
428 not further substituted in the indole ring to any extent, whether  
429 or not substituted on the cycloalkane ring to any extent \* \* \*;

430 (56) Naphthalen-1-yl  
431 1-(5-fluoropentyl)-1H-indole-3-carboxylate, also known as NM2201  
432 or CBL2201;

433 (57) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-  
434 pyrrolo[2,3-b]pyridine-3-carboxamide, also known as  
435 5F-CUMYL-P7AICA;

436 (58) 1-(4-methoxyphenyl)-N-methylpropan-2-amine, also  
437 known as para-methoxymethamphetamine or PMMA.

438 (e) **Depressants.** Unless specifically excepted or unless  
439 listed in another schedule, any material, compound, mixture, or  
440 preparation which contains any quantity of the following  
441 substances having a depressant effect on the central nervous  
442 system, including their salts, isomers, and salts of isomers,  
443 whenever the existence of such salts, isomers, and salts of  
444 isomers is possible within the specific chemical designation:

445 (1) Clonazepam,  
446 6-(2-chlorophenyl)-1-methyl-8-nitro-4H-[1,2,4]triazolo[4,3-a][1,4]  
447 benzodiazepine;

448 (2) Flualprazolam,



449 8-chloro-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4  
450 ]benzodiazepine;

451 (3) Flubromazepam,  
452 7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one;

453 (4) Flubromazolam,  
454 8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]  
455 benzodiazepin;

456 (5) Gamma-hydroxybutyric acid (other names include:  
457 GHB, gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic  
458 acid; sodium oxybate; sodium oxybutyrate);

459 (6) Mecloqualone;

460 (7) Methaqualone.

461 (f) **Stimulants.** Any material, compound, mixture or  
462 preparation which contains any quantity of the following central  
463 nervous system stimulants including optical salts, isomers and  
464 salts of isomers unless specifically excepted or unless listed in  
465 another schedule:

466 (1) Aminorex;

467 (2) N-benzylpiperazine (also known as BZP and  
468 1-benzylpiperazine);

469 (3) Cathinone;

470 (4) Fenethylamine;

471 (5) Methcathinone;

472 (6) 4-methylaminorex (also known as

473 2-amino-4-methyl-5-phenyl-2-oxazoline);



474 (7) N-ethylamphetamine;  
475 (8) Any material, compound, mixture or preparation  
476 which contains any quantity of N,N-dimethylamphetamine. (Other  
477 names include: N,N,-alpha-trimethyl-benzeneethanamine and  
478 N,N-alpha-trimethylphenethylamine);

479 (9) **Synthetic cathinones.** (A) Unless listed in  
480 another schedule, any compound other than bupropion that is  
481 structurally derived from 2-Amino-1-phenyl-1-propanone by  
482 modification in any of the following ways:

483 (i) By substitution in the phenyl ring to any  
484 extent with alkyl, alkoxy, alkylendioxy, haloalkyl or halide  
485 substituents, whether or not further substituted in the phenyl  
486 ring by one or more other univalent substituents;

487 (ii) By substitution at the 3-position with  
488 an alkyl substituent;

489 (iii) By substitution at the nitrogen atom  
490 with alkyl or dialkyl groups, or by inclusion of the nitrogen atom  
491 in a cyclic structure.

492 (B) The compounds covered in this paragraph (9)  
493 include, but are not limited to, any material, compound, mixture  
494 or preparation which contains any quantity of a synthetic  
495 cathinone found in any of the following compounds, whether or not  
496 substituted to any extent, or any of these compounds which contain  
497 any synthetic cathinone, or salts, isomers, or salts of isomers,  
498 whenever the existence of such salts, isomers or salts of isomers



499 is possible, unless specifically excepted or listed in another  
500 schedule:

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

502 (ii) 4-methyl-alpha-pyrrolidinopropiophenone  
503 ("4-MePPP");

504 (iii) Alpha-pyrrolidinopentiophenone  
505 ("α-PVP");

506 (iv)

507 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one ("butylone");

508 (v) 2-(methylamino)-1-phenylpentan-1-one  
509 ("pentedrone");

510 (vi)

511 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one

512 ("pentylone");

513 (vii) 4-fluoro-N-methylcathinone ("4-FMC");

514 (viii) 3-fluoro-N-methylcathinone ("3-FMC");

515 (ix)

516 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one ("naphyrone");

517 (x) Alpha-pyrrolidinobutiophenone ("α-PBP");

518 and

519 (xi)

520 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one

521 (N-ethylpentylone, ephylone).

522 **SECTION 2.** Section 41-29-115, Mississippi Code of 1972, is  
523 amended as follows:



524 41-29-115. (a) The controlled substances listed in this  
525 section, by whatever official name, common or usual name, chemical  
526 name, or brand name designated, are included in Schedule II.

527 **SCHEDULE II**

528 (b) **Substances, vegetable origin or chemical synthesis.**

529 Unless specifically excepted or unless listed in other schedules,  
530 any of the following substances, whether produced directly or  
531 indirectly by extraction from substances of vegetable origin, or  
532 independently by means of chemical synthesis, or by combination of  
533 extraction and chemical synthesis:

534 (1) Opium and opiate, and any salt, compound,  
535 derivative, or preparation of opium or opiate, excluding  
536 apomorphine, thebaine-derived butorphanol, dextrophan,  
537 nalbuphine, naldemedine, nalmeferone, naloxegol, naloxone and  
538 naltrexone, but including the following:

539 (i) Codeine;

540 (ii) Dihydroetorphine;

541 (iii) Ethylmorphine;

542 (iv) Etorphine hydrochloride;

543 (v) Granulated opium;

544 (vi) Hydrocodone, whether alone or in combination  
545 with any material, compound, mixture or preparation;

546 (vii) Hydromorphone;

547 (viii) Metopon;

548 (ix) Morphine;



- 549 (x) Opium extracts;
- 550 (xi) Opium fluid extracts;
- 551 (xii) Oripavine;
- 552 (xiii) Oxycodone;
- 553 (xiv) Oxymorphone;
- 554 (xv) Powdered opium;
- 555 (xvi) Raw opium;
- 556 (xvii) Thebaine;
- 557 (xviii) Tincture of opium;
- 558 (2) Any salt, compound, isomer, derivative, or
- 559 preparation thereof which is chemically equivalent or identical
- 560 with any of the substances referred to in paragraph (1), but not
- 561 including the isoquinoline alkaloids of opium;
- 562 (3) Opium poppy and poppy straw;
- 563 (4) Coca leaves and any salt, compound, derivative, or
- 564 preparation of cocaine or coca leaves, including cocaine and
- 565 ecgonine and any salt, compound, derivative, isomer, or
- 566 preparation thereof which is chemically equivalent or identical
- 567 with any of these substances, but not including:
- 568 (i) Decocainized coca leaves or extraction of coca
- 569 leaves, which extractions do not contain cocaine or ecgonine; or
- 570 (ii) Ioflupane;
- 571 (5) Concentrate of poppy straw (the crude extract of
- 572 poppy straw in either liquid, solid or powder form which contains
- 573 the phenanthrene alkaloids of the opium poppy).



574 (c) **Opiates.** Any of the following opiates, including their  
575 isomers, esters, ethers, salts, and salts of isomers, whenever the  
576 existence of these isomers, esters, ethers and salts is possible  
577 within the specified chemical designation, dextrophan and  
578 levopropoxyphene excepted:

- 579 (1) Alfentanil;
- 580 (2) Alphaprodine;
- 581 (3) Anileridine;
- 582 (4) Bezitramide;
- 583 (5) Bulk dextropropoxyphene (nondosage forms);
- 584 (6) Carfentanil;
- 585 (7) Dihydrocodeine;
- 586 (8) Diphenoxylate;
- 587 (9) Fentanyl;
- 588 (10) Isomethadone;
- 589 (11) Levo-alpha-acetylmethadol  
590 (levo-alpha-acetylmethadol, levomethadyl acetate, LAAM);
- 591 (12) Levomethorphan;
- 592 (13) Levorphanol;
- 593 (14) Metazocine;
- 594 (15) Methadone;
- 595 (16) Methadone-intermediate,  
596 4-cyano-2-dimethylamino-4,4-diphenyl butane;
- 597 (17) Moramide-intermediate,  
598 2-methyl-3-morpholino-1,1-diphenylpropane-carboxylic acid;





599 (18) Oliceridine  
600 (N-[(3-methoxythiophen-2-yl)methyl]({2-[(9R)-9-(pyridin-2-yl)-6-  
601 oxaspiro[4.5]decan-9-yl]ethyl})) amine);

602 ( \* \* \*19) Pethidine (meperidine);

603 ( \* \* \*20) Pethidine-Intermediate-A,  
604 4-cyano-1-methyl-4-phenylpiperidine;

605 ( \* \* \*21) Pethidine-Intermediate-B,  
606 ethyl-4-phenylpiperidine-4-carboxylate;

607 ( \* \* \*22) Pethidine-Intermediate-C,  
608 1-methyl-4-phenylpiperidine-4-carboxylic acid;

609 ( \* \* \*23) Phenazocine;

610 ( \* \* \*24) Piminodine;

611 ( \* \* \*25) Racemethorphan;

612 ( \* \* \*26) Racemorphan;

613 ( \* \* \*27) Remifentanil;

614 ( \* \* \*28) Sufentanil;

615 ( \* \* \*29) Tapentadol;

616 ( \* \* \*30) Thiafentanil, methyl

617 4-(2-methoxy-N-phenylacetamido)-1-(2-(thiophen-2-yl)ethyl)  
618 piperidine-4-carboxylate).

619 (d) **Stimulants.** Any material, compound, mixture, or  
620 preparation which contains any quantity of the following  
621 substances:

622 (1) Amphetamine, its salts, optical isomers, and salts  
623 of its optical isomers;



- 624 (2) Phenmetrazine and its salts;
- 625 (3) Any substance which contains any quantity of
- 626 methamphetamine, including its salts, isomers, and salts of
- 627 isomers;
- 628 (4) Methylphenidate and its salts;
- 629 (5) Lisdexamfetamine, its salts, isomers and salts of
- 630 isomers.

631 (e) **Depressants.** Unless specifically exempted or unless

632 listed in another schedule, any material, compound, mixture, or

633 preparation which contains any quantity of the following

634 substances:

- 635 (1) Amobarbital;
- 636 (2) Secobarbital;
- 637 (3) Pentobarbital;
- 638 (4) Glutethimide.

639 (f) **Hallucinogenic substances.**

- 640 (1) Dronabinol oral solution
- 641 [(-)-delta-9-trans-tetrahydrocannabinol (delta-9-THC)];
- 642 (2) Nabilone [other names include:
- 643 (+/-)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-
- 644 hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo(b,d)pyran-9-one].

645 (g) **Immediate precursors.** Unless specifically excepted or

646 unless listed in another schedule, any material, compound,

647 mixture, or preparation which contains any quantity of the

648 following substances:



649 (1) Amphetamine and methamphetamine immediate  
650 precursor: Phenylacetone (other names include:  
651 phenyl-2-propanone; P2P; benzyl methyl ketone; and methyl benzyl  
652 ketone);

653 (2) Phencyclidine immediate precursors:

654 (i) 1-phenylcyclohexylamine;

655 (ii) 1-piperidinocyclohexanecarbonitrile (PCC);

656 (3) Fentanyl immediate precursor:

657 4-anilino-N-phenethyl-4-piperidine (ANPP).

658 (h) Any material, compound, mixture or preparation which  
659 contains any quantity of a Schedule II controlled substance and is  
660 listed as an exempt substance in 21 CFR, Section 1308.24 or  
661 1308.32, shall be exempted from the provisions of the Uniform  
662 Controlled Substances Law.

663 **SECTION 3.** Section 41-29-119, Mississippi Code of 1972, is  
664 amended as follows:

665 41-29-119. (A) The controlled substances listed in this  
666 section are included in Schedule IV.

667 **SCHEDULE IV**

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



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

675 (2) Dextropropoxyphene, including its salts  
676 (Darvon, Darvon-N; also found in Darvon compound and Darvocet-N,  
677 etc.);

678 (3)  
679 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its  
680 salts, optical and geometric isomers and salts of these isomers  
681 (including tramadol).

682 (b) **Depressants.** Any material, compound, mixture or  
683 preparation which contains any quantity of the following  
684 substances:

685 (1) Alfaxalone;

686 (2) Alprazolam;

687 (3) Barbital;

688 (4) Brexanolone;

689 (5) Bromazepam;

690 (6) Camazepam;

691 (7) Carisoprodol;

692 (8) Chloral betaine;

693 (9) Chloral hydrate;

694 (10) Chlordiazepoxide and its salts, but does not  
695 include chlordiazepoxide hydrochloride and clidinium bromide or  
696 chlordiazepoxide and esterified estrogens;



- 697 (11) Clobazam;  
698 (12) Clonazepam;  
699 (13) Clorazepate;  
700 (14) Clotiazepam;  
701 (15) Cloxazolam;  
702 (16) Delorazepam;  
703 (17) Diazepam;  
704 (18) Dichloralphenazone;  
705 (19) Estazolam;  
706 (20) Ethchlorvynol;  
707 (21) Ethinamate;  
708 (22) Ethyl loflazepate;  
709 (23) Fludiazepam;  
710 (24) Flunitrazepam;  
711 (25) Flurazepam;  
712 (26) Fospropofol;  
713 (27) Halazepam;  
714 (28) Haloxazolam;  
715 (29) Ketazolam;  
716 (30) Lemborexant;  
717 ( \* \* \*31) Loprazolam;  
718 ( \* \* \*32) Lorazepam;  
719 ( \* \* \*33) Lormetazepam;  
720 ( \* \* \*34) Mebutamate;  
721 ( \* \* \*35) Medazepam;



722 ( \* \* \*36) Meprobamate;  
723 ( \* \* \*37) Methohexital;  
724 ( \* \* \*38) Methylphenobarbital;  
725 ( \* \* \*39) Midazolam;  
726 ( \* \* \*40) Nimetazepam;  
727 ( \* \* \*41) Nitrazepam;  
728 ( \* \* \*42) Nordiazepam;  
729 ( \* \* \*43) Oxazepam;  
730 ( \* \* \*44) Oxazolam;  
731 ( \* \* \*45) Paraldehyde;  
732 ( \* \* \*46) Petrichloral;  
733 ( \* \* \*47) Phenobarbital;  
734 ( \* \* \*48) Pinazepam;  
735 ( \* \* \*49) Prazepam;  
736 ( \* \* \*50) Quazepam;  
737 (51) Remimazolam;  
738 ( \* \* \*52) Suvorexant;  
739 ( \* \* \*53) Temazepam;  
740 ( \* \* \*54) Tetrazepam;  
741 ( \* \* \*55) Triazolam;  
742 ( \* \* \*56) Zaleplon;  
743 ( \* \* \*57) Zolpidem;  
744 ( \* \* \*58) Zopiclone.

745 (c) **Fenfluramine.**



746 (d) **Lorcaserin.** Any material, compound, mixture, or  
747 preparation which contains any quantity of Lorcaserin, including  
748 its salts, isomers, and salts of such isomers, whenever the  
749 existence of such salts, isomers, and salts of isomers is  
750 possible.

751 (e) **Stimulants.** Any material, compound, mixture or  
752 preparation which contains any quantity of the following  
753 substances:

- 754 (1) Cathine ((+/-) Norpseudoephedrine);
- 755 (2) Diethylpropion;
- 756 (3) Fencamfamin;
- 757 (4) Fenproporex;
- 758 (5) Mazindol;
- 759 (6) Mefenorex;
- 760 (7) Modafinil;
- 761 (8) Pemoline (including any organometallic  
762 complexes and chelates thereof);
- 763 (9) Phentermine;
- 764 (10) Pipradrol;
- 765 (11) Sibutramine;
- 766 (12) Solriamfetol;
- 767 (13) SPA ((-)-1-dimethylamino-1,2-diphenylethane).

768 (f) **Other substances.**

- 769 (1) Pentazocine;
- 770 (2) Butorphanol (including its optical isomers);



771 (3) Eluxadoline  
772 (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopr  
773 opyl][(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-meth  
774 oxybenzoic acid); (including its optical isomers) and its salts,  
775 isomers, and salts of isomers.

776 (B) Any material, compound, mixture or preparation which  
777 contains any quantity of a Schedule IV controlled substance and is  
778 listed as an exempt substance in 21 CFR, Section 1308.22, 1308.24,  
779 1308.26, 1308.32 or 1308.34, shall be exempted from the provisions  
780 of the Uniform Controlled Substances Law.

781 **SECTION 4.** Section 41-29-121, Mississippi Code of 1972, is  
782 amended as follows:

783 41-29-121.

784 **SCHEDULE V**

785 (a) Schedule V consists of the drugs and other substances,  
786 by whatever official name, common or usual name, chemical name, or  
787 brand name designated, listed in this section.

788 (b) Narcotic drugs. [Reserved]

789 (c) **Narcotic drugs containing nonnarcotic active medicinal**  
790 **ingredients.** Any compound, mixture or preparation containing any  
791 of the following narcotic drugs, or their salts calculated as the  
792 free anhydrous base or alkaloid, in limited quantities as set  
793 forth below, which also contains one or more nonnarcotic active  
794 medicinal ingredients in sufficient proportion to confer upon the





795 compound, mixture or preparation valuable medicinal qualities  
796 other than those possessed by the narcotic drug alone:

797 (1) Not more than two hundred (200) milligrams of  
798 codeine, or any of its salts, per one hundred (100) milliliters or  
799 per one hundred (100) grams;

800 (2) Not more than one hundred (100) milligrams of  
801 dihydrocodeine, or any of its salts, per one hundred (100)  
802 milliliters or per one hundred (100) grams;

803 (3) Not more than one hundred (100) milligrams of  
804 ethylmorphine, or any of its salts, per one hundred (100)  
805 milliliters or per one hundred (100) grams;

806 (4) Not more than two and five-tenths (2.5) milligrams  
807 of diphenoxylate and not less than twenty-five (25) micrograms of  
808 atropine sulphate per dosage unit;

809 (5) Not more than one hundred (100) milligrams of opium  
810 per one hundred (100) milliliters or per one hundred (100) grams;

811 (6) Not more than five-tenths (0.5) milligram of  
812 difenoxin and not less than twenty-five (25) micrograms of  
813 atropine sulfate per dosage unit.

814 (d) **Stimulants.** Unless specifically excepted or listed in  
815 another schedule, any material, compound, mixture or preparation  
816 which contains any quantity of the following substance, including  
817 its salts, isomers and salts of isomers: Pyrovalerone.

818 (e) **Depressants.** Unless specifically exempted or excluded  
819 or unless listed in another schedule, any material, compound,



820 mixture or preparation which contains any quantity of the  
821 following substances having a depressant effect on the central  
822 nervous system, including their salts, isomers and salts of  
823 isomers:

824 (1) Brivaracetam  
825 ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl] butanamide) (also  
826 referred to as BRV; UCB-34714; Briviact);

827 (2) Cenobamate  
828 [(1R)-1-(2-chlorophenyl)-2-tetrazol-2-yl]ethyl carbamate;  
829 2H-tetrazole-2-ethanol, alpha-(2-chlorophenyl)-, carbamate  
830 (ester), (alphaR)-; carbamic acid (R)-(+)-1-(2-chlorophenyl)  
831 -2-(2H-tetrazol-2-yl)ethyl ester);

832 ( \* \* \*3) Ezogabine [N-[2-amino-4-(4-  
833 fluorobenzylamino)-phenyl]-carbamic acid ethyl ester];

834 ( \* \* \*4) Lacosamide  
835 [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide];

836 (5) Lasmiditan [2,4,6-trifluoro-N-(6-  
837 (1-methylpiperidine-4-carbonyl)pyridine-2-yl-benzamide];

838 ( \* \* \*6) Pregabalin  
839 [(S)-3-(aminomethyl)-5-methylhexanoic acid].

840 (f) Any material, compound, mixture or preparation which  
841 contains any quantity of a Schedule V controlled substance and is  
842 listed as an exempt substance in 21 CFR, Section 1308.22, 1308.24,  
843 1308.26, 1308.32 or 1308.34, shall be exempted from the provisions  
844 of the Uniform Controlled Substances Law.



845           **SECTION 5.** This act shall take effect and be in force from  
846 and after July 1, 2022.

