

1                   **HOUSE OF REPRESENTATIVES - FLOOR VERSION**

2                               STATE OF OKLAHOMA

3                               1st Session of the 60th Legislature (2025)

4   ENGROSSED SENATE  
5   BILL NO. 860

By: Weaver of the Senate

and

Kannady and **Humphrey** of the  
House

10                   **[ Uniform Controlled Dangerous Substances Act -**

11                               **Schedule I substances - effective date ]**

14   BE IT ENACTED BY THE PEOPLE OF THE STATE OF OKLAHOMA:

15           SECTION 1.           AMENDATORY           63 O.S. 2021, Section 2-204, as  
16   last amended by Section 3, Chapter 308, O.S.L. 2024 (63 O.S. Supp.  
17   2024, Section 2-204), is amended to read as follows:

18           Section 2-204. The controlled substances listed in this section  
19   are included in Schedule I and include any material, compound,  
20   mixture,    or preparation that contains any quantity of the following  
21   hallucinogenic substances, their salts, isomers,    and salts of  
22   isomers, unless specifically excepted, when the existence of these  
23   salts, isomers,    and salts of isomers is possible within the specific  
24   chemical designation.

1       A. Any of the following opiates including their isomers,  
2 esters, ethers, salts, and salts of isomers, esters, and ethers,  
3 unless specifically excepted, when the existence of these isomers,  
4 esters, ethers, and salts is possible within the specific chemical  
5 designation:

- 6       1. Acetylmethadol;
- 7       2. Allylprodine;
- 8       3. Alphacetylmethadol;
- 9       4. Alphameprodine;
- 10      5. Alphamethadol;
- 11      6. Benzethidine;
- 12      7. Betacetylmethadol;
- 13      8. Betameprodine;
- 14      9. Betamethadol;
- 15      10. Betaprodine;
- 16      11. Clonitazene;
- 17      12. Dextromoramide;
- 18      13. Dextrorphan (except its methyl ether);
- 19      14. Diampromide;
- 20      15. Diethylthiambutene;
- 21      16. Dimenoxadol;
- 22      17. Dimepheptanol;
- 23      18. Dimethylthiambutene;
- 24      19. Dioxaphetyl butyrate;

1	20.	Dipipanone;
2	21.	Ethylmethylthiambutene;
3	22.	Etonitazene;
4	23.	Etoxeridine;
5	24.	Furethidine;
6	25.	Hydroxypethidine;
7	26.	Isotonitazene;
8	27.	Ketobemidone;
9	28.	Levomoramide;
10	29.	Levophenacymorphan;
11	30.	Metonitazene;
12	31.	Morpheridine;
13	32.	N-desethyl isotonitazene;
14	33.	N-pyrrolidino protonitazene;
15	34.	Noracymethadol;
16	35.	Norlevorphanol;
17	36.	Normethadone;
18	37.	Norpipanone;
19	38.	Phenadoxone;
20	39.	Phenampromide;
21	40.	Phenomorphane;
22	41.	Phenoperidine;
23	42.	Piritramide;
24	43.	Proheptazine;

1 44. Properidine;

2 45. Protonitazene;

3 46. Racemoramide; or

4 47. Trimeperidine.

5 B. Any of the following opium derivatives, their salts,  
6 isomers, and salts of isomers, unless specifically excepted, when  
7 the existence of these salts, isomers, and salts of isomers is  
8 possible within the specific chemical designation:

9 1. Acetorphine;

10 2. Acetyldihydrocodeine;

11 3. Benzylmorphine;

12 4. Codeine methylbromide;

13 5. Codeine-N-Oxide;

14 6. Cyprenorphine;

15 7. Desomorphine;

16 8. Dihydromorphine;

17 9. Etorphine;

18 10. Heroin;

19 11. Hydromorphinol;

20 12. Methyldesorphine;

21 13. Methylhydromorphine;

22 14. Morphine methylbromide;

23 15. Morphine methylsulfonate;

24 16. Morphine-N-Oxide;

- 1 17. Myrophine;
- 2 18. Nicocodeine;
- 3 19. Nicomorphine;
- 4 20. Normorphine;
- 5 21. Phoclodine;
- 6 22. Thebacon;
- 7 23. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide
- 8 (Acetyl fentanyl);
- 9 24. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butenamide
- 10 (Crotonyl fentanyl);
- 11 25. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-
- 12 furancarboxamide (Furanyl fentanyl);
- 13 26. N-phenyl-1-(2-phenylethyl)-4-piperidinamine (4-ANPP);
- 14 27. N-(1-phenethylpiperidin-4-yl)-N-
- 15 phenylcyclopropanecarboxamide (Cyclopropyl fentanyl); or
- 16 28. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide
- 17 (Butyrl fentanyl).
- 18 C. Any material, compound, mixture, or preparation which
- 19 contains any quantity of the following hallucinogenic substances,
- 20 their salts, isomers, and salts of isomers, unless specifically
- 21 excepted, when the existence of these salts, isomers, and salts of
- 22 isomers is possible within the specific chemical designation:
- 23 1. Methcathinone;
- 24 2. 3, 4-methylenedioxy amphetamine;

- 1 3. 3, 4-methylenedioxy methamphetamine;
- 2 4. 5-methoxy-3, 4-methylenedioxy amphetamine;
- 3 5. 3, 4, 5-trimethoxy amphetamine;
- 4 6. Bufotenine;
- 5 7. Diethyltryptamine;
- 6 8. Dimethyltryptamine;
- 7 9. 4-methyl-2, 5-dimethoxyamphetamine;
- 8 10. Ibogaine;
- 9 11. Lysergic acid diethylamide;
- 10 12. Marijuana;
- 11 13. Mescaline;
- 12 14. N-benzylpiperazine;
- 13 15. N-ethyl-3-piperidyl benzilate;
- 14 16. N-methyl-3-piperidyl benzilate;
- 15 17. Psilocybin;
- 16 18. Psilocyn;
- 17 19. 2, 5 dimethoxyamphetamine;
- 18 20. 4 Bromo-2, 5-dimethoxyamphetamine;
- 19 21. 4 methoxyamphetamine;
- 20 22. Cyclohexamine;
- 21 23. Salvia Divinorum;
- 22 24. Salvinorin A;

- 1        25. Thiophene Analog of Phencyclidine. Also known as: 1-(1-(2-  
2 thienyl) cyclohexyl) piperidine; 2-Thienyl Analog of Phencyclidine;  
3 TCP, TCP;  
4        26. Phencyclidine (PCP);  
5        27. Pyrrolidine Analog for Phencyclidine. Also known as 1-(1-  
6 Phenylcyclohexyl) - Pyrrolidine, PCPy, PHP;  
7        28. 1-(3-trifluoromethylphenyl) piperazine;  
8        29. Flunitrazepam;  
9        30. B-hydroxy-amphetamine;  
10       31. B-ketoamphetamine;  
11       32. 2,5-dimethoxy-4-nitroamphetamine;  
12       33. 2,5-dimethoxy-4-bromophenethylamine;  
13       34. 2,5-dimethoxy-4-chlorophenethylamine;  
14       35. 2,5-dimethoxy-4-iodoamphetamine;  
15       36. 2,5-dimethoxy-4-iodophenethylamine;  
16       37. 2,5-dimethoxy-4-methylphenethylamine;  
17       38. 2,5-dimethoxy-4-ethylphenethylamine;  
18       39. 2,5-dimethoxy-4-fluorophenethylamine;  
19       40. 2,5-dimethoxy-4-nitrophenethylamine;  
20       41. 2,5-dimethoxy-4-ethylthio-phenethylamine;  
21       42. 2,5-dimethoxy-4-isopropylthio-phenethylamine;  
22       43. 2,5-dimethoxy-4-propylthio-phenethylamine;  
23       44. 2,5-dimethoxy-4-cyclopropylmethylthio-phenethylamine;  
24       45. 2,5-dimethoxy-4-tert-butylthio-phenethylamine;

- 1 46. 2,5-dimethoxy-4-(2-fluoroethylthio)-phenethylamine;
- 2 47. 5-methoxy-N, N-dimethyltryptamine;
- 3 48. N-methyltryptamine;
- 4 49. A-ethyltryptamine;
- 5 50. A-methyltryptamine;
- 6 51. N, N-diethyltryptamine;
- 7 52. N, N-diisopropyltryptamine;
- 8 53. N, N-dipropyltryptamine;
- 9 54. 5-methoxy-a-methyltryptamine;
- 10 55. 4-hydroxy-N, N-diethyltryptamine;
- 11 56. 4-hydroxy-N, N-diisopropyltryptamine;
- 12 57. 5-methoxy-N, N-diisopropyltryptamine;
- 13 58. 4-hydroxy-N-isopropyl-N-methyltryptamine;
- 14 59. 3,4-Methylenedioxymethcathinone (Methylone);
- 15 60. 3,4-Methylenedioxypyrovalerone (MDPV);
- 16 61. 3-Methylmethcathinone (Metaphedrone);
- 17 62. 4-Methylmethcathinone (Mephedrone);
- 18 63. 4-methoxymethcathinone;
- 19 64. 4-Fluoromethcathinone;
- 20 65. 3-Fluoromethcathinone;
- 21 66. 1-(8-bromobenzo 1,2-b;4,5-b' difuran-4-yl)-2-aminopropane;
- 22 67. 2,5-Dimethoxy-4-chloroamphetamine;
- 23 68. 4-Methylethcathinone;
- 24 69. Pyrovalerone;



1 70. N,N-diallyl-5-methoxytryptamine;  
 2 71. 3,4-Methylenedioxy-N-ethylcathinone (Ethylone);  
 3 72. B-keto-N-Methylbenzodioxolylbutanamine (Butylone);  
 4 73. B-keto-Methylbenzodioxolylpentanamine (Pentylone);  
 5 74. Alpha-Pyrrolidinopentiophenone;  
 6 75. 4-Fluoroamphetamine;  
 7 76. Pentedrone;  
 8 77. 4'-Methyl-a-pyrrolidinohexaphenone;  
 9 78. 2,5-dimethoxy-4-(n)-propylphenethylamine;  
 10 79. 2,5-dimethoxyphenethylamine;  
 11 80. 1,4-Dibenzylpiperazine;  
 12 81. N,N-Dimethylamphetamine;  
 13 82. 4-Fluoromethamphetamine;  
 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine  
 15 (25C-NBOMe);  
 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine  
 17 (25I-NBOMe);  
 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine  
 19 (25B-NBOMe);  
 20 86. 1-(4-Fluorophenyl)piperazine;  
 21 87. Methoxetamine;  
 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N-  
 23 methylbenzamide;  
 24 89. N-ethyl hexadrone;

90. Isopropyl-U-47700;
91. Para-fluorobutyl fentanyl;
92. Para-fluorofentanyl (pFF);
93. Fluoro isobutyl fentanyl;
94. 3-Hydroxy Phencyclidine (PCP);
95. 3-methoxy Phencyclidine (PCP);
96. Flualprazolam; or
97. Flubromazolam.

D. Unless specifically excepted or unless listed in a different schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having stimulant or depressant effect on the central nervous system:

1. Fenethylline;
2. Mecloqualone;
3. N-ethylamphetamine;
4. Methaqualone;
5. Gamma-Hydroxybutyric Acid, also known as GHB, gamma-hydroxybutyrate, 4-hydroxybutyrate, 4-hydroxybutanoic acid, sodium oxybate, and sodium oxybutyrate;
6. Gamma-Butyrolactone (GBL) as packaged, marketed, manufactured, or promoted for human consumption, with the exception of legitimate food additive and manufacturing purposes;

1        7. Gamma Hydroxyvalerate (GHV) as packaged, marketed, or  
2 manufactured for human consumption, with the exception of legitimate  
3 food additive and manufacturing purposes;

4        8. Gamma Valerolactone (GVL) as packaged, marketed, or  
5 manufactured for human consumption, with the exception of legitimate  
6 food additive and manufacturing purposes;

7        9. 1,4 Butanediol (1,4 BD or BDO) as packaged, marketed,  
8 manufactured, or promoted for human consumption with the exception  
9 of legitimate manufacturing purposes; or

10       10. N-ethylpentylone.

11       E. 1. The following industrial uses of Gamma-Butyrolactone,  
12 Gamma Hydroxyvalerate, Gamma Valerolactone, or 1,4 Butanediol are  
13 excluded from all schedules of controlled substances under this  
14 title:

- 15           a. pesticides,
- 16           b. photochemical etching,
- 17           c. electrolytes of small batteries or capacitors,
- 18           d. viscosity modifiers in polyurethane,
- 19           e. surface etching of metal coated plastics,
- 20           f. organic paint disbursements for water soluble inks,
- 21           g. pH regulators in the dyeing of wool and polyamide  
22           fibers,
- 23           h. foundry chemistry as a catalyst during curing,
- 24

- i. curing agents in many coating systems based on urethanes and amides,
- j. additives and flavoring agents in food, confectionary, and beverage products,
- k. synthetic fiber and clothing production,
- l. tetrahydrofuran production,
- m. gamma butyrolactone production,
- n. polybutylene terephthalate resin production,
- o. polyester raw materials for polyurethane elastomers and foams,
- p. coating resin raw material, and
- q. as an intermediate in the manufacture of other chemicals and pharmaceuticals.

2. At the request of any person, the Director of the Oklahoma State Bureau of Narcotics and Dangerous Drugs Control may exempt any other product containing Gamma-Butyrolactone, Gamma Hydroxyvalerate, Gamma Valerolactone, or 1,4 Butanediol from being included as a Schedule I controlled substance if such product is labeled, marketed, manufactured, and distributed for legitimate industrial use in a manner that reduces or eliminates the likelihood of abuse.

3. In making a determination regarding an industrial product, the Director, after notice and hearing, shall consider the following:

- a. the history and current pattern of abuse,

- b. the name and labeling of the product,
- c. the intended manner of distribution, advertising, and promotion of the product, and
- d. other factors as may be relevant to and consistent with the public health and safety.

4. The hearing shall be held in accordance with the procedures of the Administrative Procedures Act.

F. Any material, compound, mixture, or preparation, whether produced directly or indirectly from a substance of vegetable origin or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis, that contains any quantity of the following substances, or that contains any of their salts, isomers, and salts of isomers when the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

1. JWH-004;
2. JWH-007;
3. JWH-009;
4. JWH-015;
5. JWH-016;
6. JWH-018;
7. JWH-019;
8. JWH-020;
9. JWH-030;

1	10.	JWH-046;
2	11.	JWH-047;
3	12.	JWH-048;
4	13.	JWH-049;
5	14.	JWH-050;
6	15.	JWH-070;
7	16.	JWH-071;
8	17.	JWH-072;
9	18.	JWH-073;
10	19.	JWH-076;
11	20.	JWH-079;
12	21.	JWH-080;
13	22.	JWH-081;
14	23.	JWH-082;
15	24.	JWH-094;
16	25.	JWH-096;
17	26.	JWH-098;
18	27.	JWH-116;
19	28.	JWH-120;
20	29.	JWH-122;
21	30.	JWH-145;
22	31.	JWH-146;
23	32.	JWH-147;
24	33.	JWH-148;

1	34.	JWH-149;
2	35.	JWH-150;
3	36.	JWH-156;
4	37.	JWH-167;
5	38.	JWH-175;
6	39.	JWH-180;
7	40.	JWH-181;
8	41.	JWH-182;
9	42.	JWH-184;
10	43.	JWH-185;
11	44.	JWH-189;
12	45.	JWH-192;
13	46.	JWH-193;
14	47.	JWH-194;
15	48.	JWH-195;
16	49.	JWH-196;
17	50.	JWH-197;
18	51.	JWH-198;
19	52.	JWH-199;
20	53.	JWH-200;
21	54.	JWH-201;
22	55.	JWH-202;
23	56.	JWH-203;
24	57.	JWH-204;

1	58.	JWH-205;
2	59.	JWH-206;
3	60.	JWH-207;
4	61.	JWH-208;
5	62.	JWH-209;
6	63.	JWH-210;
7	64.	JWH-211;
8	65.	JWH-212;
9	66.	JWH-213;
10	67.	JWH-234;
11	68.	JWH-235;
12	69.	JWH-236;
13	70.	JWH-237;
14	71.	JWH-239;
15	72.	JWH-240;
16	73.	JWH-241;
17	74.	JWH-242;
18	75.	JWH-243;
19	76.	JWH-244;
20	77.	JWH-245;
21	78.	JWH-246;
22	79.	JWH-248;
23	80.	JWH-249;
24	81.	JWH-250;



1	82.	JWH-251;
2	83.	JWH-252;
3	84.	JWH-253;
4	85.	JWH-262;
5	86.	JWH-292;
6	87.	JWH-293;
7	88.	JWH-302;
8	89.	JWH-303;
9	90.	JWH-304;
10	91.	JWH-305;
11	92.	JWH-306;
12	93.	JWH-307;
13	94.	JWH-308;
14	95.	JWH-311;
15	96.	JWH-312;
16	97.	JWH-313;
17	98.	JWH-314;
18	99.	JWH-315;
19	100.	JWH-316;
20	101.	JWH-346;
21	102.	JWH-348;
22	103.	JWH-363;
23	104.	JWH-364;
24	105.	JWH-365;

1	106.	JWH-367;
2	107.	JWH-368;
3	108.	JWH-369;
4	109.	JWH-370;
5	110.	JWH-371;
6	111.	JWH-373;
7	112.	JWH-386;
8	113.	JWH-387;
9	114.	JWH-392;
10	115.	JWH-394;
11	116.	JWH-395;
12	117.	JWH-397;
13	118.	JWH-398;
14	119.	JWH-399;
15	120.	JWH-400;
16	121.	JWH-412;
17	122.	JWH-413;
18	123.	JWH-414;
19	124.	JWH-415;
20	125.	CP-55, 940;
21	126.	CP-47, 497;
22	127.	HU-210;
23	128.	HU-211;
24	129.	WIN-55, 212-2;

1 130. AM-2201;  
2 131. AM-2233;  
3 132. JWH-018 adamantyl-carboxamide;  
4 133. AKB48;  
5 134. JWH-122 N-(4-pentenyl) analog;  
6 135. MAM2201;  
7 136. URB597;  
8 137. URB602;  
9 138. URB754;  
10 139. UR144;  
11 140. XLR11;  
12 141. A-796,260;  
13 142. STS-135;  
14 143. AB-FUBINACA;  
15 144. AB-PINACA;  
16 145. PB-22;  
17 146. AKB48 N-5-Fluoropentyl;  
18 147. AM1248;  
19 148. FUB-PB-22;  
20 149. ADB-FUBINACA;  
21 150. BB-22;  
22 151. 5-Fluoro PB-22; or  
23 152. 5-Fluoro AKB-48.  
24

1 G. In addition to those substances listed in subsection F of  
2 this section, unless specifically excepted or unless listed in  
3 another schedule, any material, compound, mixture, or preparation  
4 which contains any quantity of a synthetic cannabinoid found to be  
5 in any of the following chemical groups:

6 1. Naphthoylindoles: any compound containing a 3-(1-  
7 naphthoyl)indole structure with or without substitution at the  
8 nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,  
9 alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-  
10 (N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-  
11 2-pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl,  
12 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or  
13 halophenyl group, whether or not further substituted on the indole  
14 ring to any extent, and whether or not substituted on the naphthyl  
15 ring to any extent. Naphthoylindoles include, but are not limited  
16 to:

- 17 a. 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWH-  
18 200),
- 19 b. 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201),
- 20 c. 1-pentyl-3-(1-naphthoyl)indole (JWH-018),
- 21 d. 1-butyl-3-(1-naphthoyl)indole (JWH-073),
- 22 e. 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (JWH-081),
- 23 f. 1-propyl-2-methyl-3-(1-naphthoyl)indole (JWH-015),
- 24 g. 1-hexyl-3-(1-naphthoyl)indole (JWH-019),

- h. 1-pentyl-3-(4-methyl-1-naphthoyl)indole (JWH-122),  
i. 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (JWH-210),  
j. 1-pentyl-3-(4-chloro-1-naphthoyl)indole (JWH-398),  
k. 1-pentyl-2-methyl-3-(1-naphthoyl)indole (JWH-007),  
l. 1-pentyl-3-(7-methoxy-1-naphthoyl)indole (JWH-164),  
m. 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole  
(JWH-098),  
n. 1-pentyl-3-(4-fluoro-1-naphthoyl)indole (JWH-412),  
o. 1-[1-(N-methyl-2-piperidinyl)methyl]-3-(1-naphthoyl)indole (AM-1220),  
p. 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole  
(MAM-2201), or  
q. 1-(4-cyanobutyl)-3-(1-naphthoyl)indole (AM-2232);

2. Naphthylmethylindeles: any compound containing a 1H-indol-3-yl-(1-naphthyl)methane structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not further substituted on the indole ring to any extent, and whether or not substituted on the naphthyl ring to any extent. Naphthylmethylindeles include, but are not limited to, (1-pentylindol-3-yl)(1-naphthyl)methane (JWH-175);

1        3. Naphthoylpyrroles: any compound containing a 3-(1-  
2 naphthoyl)pyrrole structure with or without substitution at the  
3 nitrogen atom of the pyrrole ring by an alkyl, haloalkyl,  
4 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,  
5 halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-  
6 morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
7 morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl,  
8 phenyl, or halophenyl group, whether or not further substituted on  
9 the pyrrole ring to any extent, and whether or not substituted on  
10 the naphthyl group to any extent. Naphthoylpyrroles include, but  
11 are not limited to:

- 12            a. 1-hexyl-2-phenyl-4-(1-naphthoyl)pyrrole (JWH-147),
- 13            b. 1-pentyl-5-(2-methylphenyl)-3-(1-naphthoyl)pyrrole  
14                    (JWH-370),
- 15            c. 1-pentyl-3-(1-naphthoyl)pyrrole (JWH-030), or
- 16            d. 1-hexyl-5-phenyl-3-(1-naphthoyl)pyrrole (JWH-147);

17        4. Naphthylideneindenes: any compound containing a 1-(1-  
18 naphthylmethylene)indene structure with or without substitution at  
19 the 3-position of the indene ring by an alkyl, haloalkyl,  
20 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,  
21 halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-  
22 morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
23 morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl,  
24 phenyl, or halophenyl group, whether or not further substituted on

1 the indene group to any extent, and whether or not substituted on  
2 the naphthyl group to any extent. Naphthylmethylindenes include,  
3 but are not limited to, (1-[(3-pentyl)-1H-inden-1-  
4 ylidene)methyl]naphthalene (JWH-176);

5 5. Phenylacetylindoles: any compound containing a 3-  
6 phenylacetylindole structure with or without substitution at the  
7 nitrogen atom of the indole ring by alkyl, haloalkyl, cyanoalkyl,  
8 alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-  
9 (N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-  
10 2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl,  
11 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or  
12 halophenyl group, whether or not further substituted on the indole  
13 ring to any extent, and whether or not substituted on the phenyl  
14 ring to any extent. Phenylacetylindoles include, but are not  
15 limited to:

- 16 a. 1-pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250),
- 17 b. 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole  
18 (RCS-8),
- 19 c. 1-pentyl-3-(2-chlorophenylacetyl)indole (JWH-203),
- 20 d. 1-pentyl-3-(2-methylphenylacetyl)indole (JWH-251),
- 21 e. 1-pentyl-3-(4-methoxyphenylacetyl)indole (JWH-201), or
- 22 f. 1-pentyl-3-(3-methoxyphenylacetyl)indole (JWH-302);

23 6. Cyclohexylphenols: any compound containing a 2-(3-  
24 hydroxycyclohexyl)phenol structure with or without substitution at

the 5-position of the phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, and whether or not further substituted on the cyclohexyl ring to any extent. Cyclohexylphenols include, but are not limited to:

- a. 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol (CP-47,497),
- b. 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol (cannabicyclohexanol; CP-47,497 C8 homologue),  
or
- c. 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol (CP 55, 940);

7. Benzoylindoles: any compound containing a 3-(benzoyl)indole structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not further substituted on the indole ring to any extent, and whether or not substituted on the phenyl



group to any extent. Benzoylindoles include, but are not limited to:

- a. 1-pentyl-3-(4-methoxybenzoyl)indole (RCS-4),
- b. 1-[2-(4-morpholinyl)ethyl]-2-methyl-3-(4-methoxybenzoyl)indole (Pravadoline or WIN 48, 098),
- c. 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM-694),
- d. 1-pentyl-3-(2-iodobenzoyl)indole (AM-679), or
- e. 1-[1-(N-methyl-2-piperidinyl)methyl]-3-(2-iodobenzoyl)indole (AM-2233);

8. Cyclopropoylindoles: Any compound containing a 3-(cyclopropoyl)indole structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not further substituted in the indole ring to any extent, and whether or not substituted in the cyclopropoyl ring to any extent. Cyclopropoylindoles include, but are not limited to:

- a. 1-pentyl-3-(2,2,3,3-tetramethylcyclopropoyl)indole (UR-144),
- b. 1-(5-chloropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole (5Cl-UR-144), or

c. 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole (XLR11);

9. Indole Amides: Any compound containing a 1H-Indole-3-carboxamide structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not substituted at the carboxamide group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl, cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not further substituted in the indole, adamantyl, naphthyl, phenyl, pyrrole, quinolinyl, or cycloalkyl rings to any extent. Indole Amides include, but are not limited to:

a. N-(1-adamantyl)-1-pentyl-1H-indole-3-carboxamide (2NE1),

b. N-(1-adamantyl)-1-(5-fluoropentyl)-1H-indole-3-carboxamide (STS-135),

c. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-carboxamide (ADBICA),

- d. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indole-3-carboxamide (5F-ADBICA),
- e. N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide (NNE1),
- f. 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide (5F-NNE1),
- g. N-benzyl-1-pentyl-1H-indole-3-carboxamide (SDB-006),  
or
- h. N-benzyl-1-(5-fluoropentyl)-1H-indole-3-carboxamide (5F-SDB-006);

10. Indole Esters: Any compound containing a 1H-Indole-3-carboxylate structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not substituted at the carboxylate group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl, cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not further substituted in the indole, adamantyl, naphthyl, phenyl,

pyrrole, quinolinyl, or cycloalkyl rings to any extent. Indole Esters include, but are not limited to:

- a. quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB-22),
- b. quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5F-PB-22),
- c. quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate (BB-22),
- d. naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate (FDU-PB-22), or
- e. naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (NM2201);

11. Adamantanoylindoles: Any compound containing an adamantanyl-(1H-indol-3-yl)methanone structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not further substituted in the indole ring to any extent, and whether or not substituted in the adamantyl ring to any extent. Adamantanoylindoles include, but are not limited to:

- 1           a.    adamantan-1-yl[1-[(1-methyl-2-piperidinyl)methyl]-1H-  
2               indol-3-yl]methanone (AM1248), or  
3           b.    adamantan-1-yl-(1-pentyl-1H-indol-3-yl)methanone (AB-  
4               001);

5       12.   Carbazole Ketone: Any compound containing (9H-carbazole-3-  
6   yl) methanone structure with or without substitution at the nitrogen  
7   atom of the carbazole ring by an alkyl, haloalkyl, cyanoalkyl,  
8   alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-  
9   (N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-  
10 2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl,  
11 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or  
12 halophenyl group, with substitution at the carbon of the methanone  
13 group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,  
14 cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-  
15 1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-  
16 dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not  
17 further substituted at the carbazole, adamantyl, naphthyl, phenyl,  
18 pyrrole, quinolinyl, or cycloalkyl rings to any extent. Carbazole  
19 Ketones include, but are not limited to, naphthalen-1-yl(9-pentyl-  
20 9H-carbazol-3-yl)methanone (EG-018);

21       13.   Benzimidazole Ketone: Any compound containing  
22   (benzimidazole-2-yl) methanone structure with or without  
23   substitution at either nitrogen atom of the benzimidazole ring by an  
24   alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,

1 cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-  
2 piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-  
3 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl,  
4 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or  
5 halophenyl group, with substitution at the carbon of the methanone  
6 group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,  
7 cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-  
8 1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-  
9 dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not  
10 further substituted in the benzimidazole, adamantyl, naphthyl,  
11 phenyl, pyrrole, quinolinyl, or cycloalkyl rings to any extent.

12 Benzimidazole Ketones include, but are not limited to:

- 13 a. naphthalen-1-yl(1-pentyl-1H-benzo[d]imidazol-2-  
14 1)methanone (JWH-018 benzimidazole analog), or
- 15 b. (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-  
16 yl)(naphthalen-1-yl)methanone (FUBIMINA); and

17 14. Modified by Replacement: any compound defined in this  
18 subsection that is modified by replacement of a carbon with nitrogen  
19 in the indole, naphthyl, indene, benzimidazole, or carbazole ring.

20 H. Any material, compound, mixture, extract, or preparation  
21 that contains a prohibited kratom product as provided in paragraphs  
22 3 and 4 of subsection A of Section 1-1432.4 of this title.

23 I. Any prescription drug approved by the federal Food and Drug  
24 Administration under the provisions of Section 505 of the Federal

1 Food, Drug, and Cosmetic Act, Title 21 of the United States Code,  
2 Section 355, that is designated, rescheduled, or deleted as a  
3 controlled substance under federal law by the United States Drug  
4 Enforcement Administration shall be excluded from Schedule I and  
5 shall be prescribed, distributed, dispensed, or used in accordance  
6 with federal law upon the issuance of a notice, final rule, or  
7 interim final rule by the United States Drug Enforcement  
8 Administration designating, rescheduling, or deleting as a  
9 controlled substance such a drug product under federal law, unless  
10 and until the State Board of Pharmacy takes action pursuant to  
11 Section 2-201 of this title. If the Board of Pharmacy does not take  
12 action pursuant to Section 2-201 of this title, the drug product  
13 shall be deemed to be designated, rescheduled, or deleted as a  
14 controlled substance in accordance with federal law and in  
15 compliance with the Uniform Controlled Dangerous Substances Act.

16 SECTION 2. This act shall become effective November 1, 2025.

17  
18 COMMITTEE REPORT BY: COMMITTEE ON JUDICIARY AND PUBLIC SAFETY  
19 OVERSIGHT, dated 04/22/2025 - DO PASS, As Amended and Coauthored.  
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