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AS AMENDED

By: Standridge

[Uniform Controlled Dangerous Substances Act -
Schedule I - Schedule IV - effective date]

SECTION 1. AMENDATORY 63 O.S. 2011, Section 2-204, as last amended by Section 1, Chapter 207, O.S.L. 2019 (63 O.S. Supp. 2019, Section 2-204), is amended to read as follows:

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

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

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

25. Hydroxypethidine;
26. Ketobemidone;
27. Levomoramide;
28. Levophenacymorphan;
29. Morpheridine;
30. Noracymethadol;
31. Norlevorphanol;
32. Normethadone;
33. Norpipanone;
34. Phenadoxone;
35. Phenampromide;
36. Phenomorphan;
37. Phenoperidine;
38. Piritramide;
39. Proheptazine;
40. Properidine;
41. Racemoramide; or
42. Trimeperidine.

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

1. Acetorphine;
2. Acetyldihydrocodeine;

- 1 3. Benzylmorphine;
- 2 4. Codeine methylbromide;
- 3 5. Codeine-N-Oxide;
- 4 6. Cyprenorphine;
- 5 7. Desomorphine;
- 6 8. Dihydromorphine;
- 7 9. Etorphine;
- 8 10. Heroin;
- 9 11. Hydromorphenol;
- 10 12. Methyldesorphine;
- 11 13. Methylhydromorphine;
- 12 14. Morphine methylbromide;
- 13 15. Morphine methylsulfonate;
- 14 16. Morphine-N-Oxide;
- 15 17. Myrophine;
- 16 18. Nicocodeine;
- 17 19. Nicomorphine;
- 18 20. Normorphine;
- 19 21. Phoclodine;
- 20 22. Thebacon;
- 21 23. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide
- 22 (Acetyl fentanyl);
- 23 24. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butenamide
- 24 (Crotonyl fentanyl);

1 25. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-
2 furancarboxamide (Furanyl fentanyl);
3 26. N-phenyl-1-(2-phenylethyl)-4-piperidinamine (4-ANPP);
4 27. N-(1-phenethylpiperidin-4-yl)-N-
5 phenylcyclopropanecarboxamide (Cyclopropyl fentanyl); or
6 28. N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide
7 (Butyl fentanyl).

8 C. Any material, compound, mixture, or preparation which
9 contains any quantity of the following hallucinogenic substances,
10 their salts, isomers, and salts of isomers, unless specifically
11 excepted, when the existence of these salts, isomers, and salts of
12 isomers is possible within the specific chemical designation:

- 13 1. Methcathinone;
- 14 2. 3, 4-methylenedioxy amphetamine;
- 15 3. 3, 4-methylenedioxy methamphetamine;
- 16 4. 5-methoxy-3, 4-methylenedioxy amphetamine;
- 17 5. 3, 4, 5-trimethoxy amphetamine;
- 18 6. Bufotenine;
- 19 7. Diethyltryptamine;
- 20 8. Dimethyltryptamine;
- 21 9. 4-methyl-2, 5-dimethoxyamphetamine;
- 22 10. Ibogaine;
- 23 11. Lysergic acid diethylamide;
- 24 12. Marihuana;

- 1 13. Mescaline;
- 2 14. N-benzylpiperazine;
- 3 15. N-ethyl-3-piperidyl benzilate;
- 4 16. N-methyl-3-piperidyl benzilate;
- 5 17. Psilocybin;
- 6 18. Psilocyn;
- 7 19. 2, 5 dimethoxyamphetamine;
- 8 20. 4 Bromo-2, 5-dimethoxyamphetamine;
- 9 21. 4 methoxyamphetamine;
- 10 22. Cyclohexamine;
- 11 23. Salvia Divinorum;
- 12 24. Salvinorin A;
- 13 25. Thiophene Analog of Phencyclidine. Also known as: 1-(1-(2-
- 14 thienyl) cyclohexyl) piperidine; 2-Thienyl Analog of Phencyclidine;
- 15 TCP, TCP, Tenocyclidine;
- 16 26. Phencyclidine (PCP);
- 17 27. Pyrrolidine Analog for Phencyclidine. Also known as 1-(1-
- 18 Phenylcyclohexyl) - Pyrrolidine, PCPy, PHP;
- 19 28. 1-(3-trifluoromethylphenyl) piperazine;
- 20 29. Flunitrazepam;
- 21 30. B-hydroxy-amphetamine;
- 22 31. B-ketoamphetamine;
- 23 32. 2,5-dimethoxy-4-nitroamphetamine;
- 24 33. 2,5-dimethoxy-4-bromophenethylamine;

- 1 34. 2,5-dimethoxy-4-chlorophenethylamine;
- 2 35. 2,5-dimethoxy-4-iodoamphetamine;
- 3 36. 2,5-dimethoxy-4-iodophenethylamine;
- 4 37. 2,5-dimethoxy-4-methylphenethylamine;
- 5 38. 2,5-dimethoxy-4-ethylphenethylamine;
- 6 39. 2,5-dimethoxy-4-fluorophenethylamine;
- 7 40. 2,5-dimethoxy-4-nitrophenethylamine;
- 8 41. 2,5-dimethoxy-4-ethylthio-phenethylamine;
- 9 42. 2,5-dimethoxy-4-isopropylthio-phenethylamine;
- 10 43. 2,5-dimethoxy-4-propylthio-phenethylamine;
- 11 44. 2,5-dimethoxy-4-cyclopropylmethylthio-phenethylamine;
- 12 45. 2,5-dimethoxy-4-tert-butylthio-phenethylamine;
- 13 46. 2,5-dimethoxy-4-(2-fluoroethylthio)-phenethylamine;
- 14 47. 5-methoxy-N, N-dimethyltryptamine;
- 15 48. N-methyltryptamine;
- 16 49. A-ethyltryptamine;
- 17 50. A-methyltryptamine;
- 18 51. N, N-diethyltryptamine;
- 19 52. N, N-diisopropyltryptamine;
- 20 53. N, N-dipropyltryptamine;
- 21 54. 5-methoxy-a-methyltryptamine;
- 22 55. 4-hydroxy-N, N-diethyltryptamine;
- 23 56. 4-hydroxy-N, N-diisopropyltryptamine;
- 24 57. 5-methoxy-N, N-diisopropyltryptamine;

- 1 58. 4-hydroxy-N-isopropyl-N-methyltryptamine;
- 2 59. 3,4-Methylenedioxyethcathinone (Methylone);
- 3 60. 3,4-Methylenedioxypropylvalerone (MDPV);
- 4 61. 4-Methylmethcathinone (Mephedrone);
- 5 62. 4-methoxymethcathinone;
- 6 63. 4-Fluoromethcathinone;
- 7 64. 3-Fluoromethcathinone;
- 8 65. 1-(8-bromobenzo 1,2-b;4,5-b' difuran-4-yl)-2-aminopropane;
- 9 66. 2,5-Dimethoxy-4-chloroamphetamine;
- 10 67. 4-Methylethcathinone;
- 11 68. Pyrovalerone;
- 12 69. N,N-diallyl-5-methoxytryptamine;
- 13 70. 3,4-Methylenedioxy-N-ethylcathinone (Ethylone);
- 14 71. B-keto-N-Methylbenzodioxolylbutanamine (Butylone);
- 15 72. B-keto-Methylbenzodioxolylpentanamine (Pentylone);
- 16 73. Alpha-Pyrrolidinopentiophenone;
- 17 74. 4-Fluoroamphetamine;
- 18 75. Pentedrone;
- 19 76. 4'-Methyl-a-pyrrolidinohexaphenone;
- 20 77. 2,5-dimethoxy-4-(n)-propylphenethylamine;
- 21 78. 2,5-dimethoxyphenethylamine;
- 22 79. 1,4-Dibenzylpiperazine;
- 23 80. N,N-Dimethylamphetamine;
- 24 81. 4-Fluoromethamphetamine;

- 1 82. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine
2 (25C-NBOMe) ;
- 3 83. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine
4 (25I-NBOMe) ;
- 5 84. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine
6 (25B-NBOMe) ;
- 7 85. 1-(4-Fluorophenyl)piperazine;
- 8 86. Methoxetamine;
- 9 87. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N-
10 methylbenzamide;
- 11 88. N-ethyl hexadrone;
- 12 89. Isopropyl-U-47700;
- 13 90. Para-fluorobutyl fentanyl;
- 14 91. Fluoro isobutryl fentanyl;
- 15 92. 3-Hydroxy Phencyclidine (PCP); ~~or~~
- 16 93. 3-methoxy Phencyclidine (PCP); i
- 17 94. Flualprazolam; or
- 18 95. Flubromazolam.

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

- 23 1. Fenethylamine;
- 24 2. Mecloqualone;

1 3. N-ethylamphetamine;

2 4. Methaqualone;

3 5. Gamma-Hydroxybutyric Acid, also known as GHB, gamma-
4 hydroxybutyrate, 4-hydroxybutyrate, 4-hydroxybutanoic acid, sodium
5 oxybate~~7~~ and sodium oxybutyrate;

6 6. Gamma-Butyrolactone (GBL) as packaged, marketed,
7 manufactured or promoted for human consumption, with the exception
8 of legitimate food additive and manufacturing purposes;

9 7. Gamma Hydroxyvalerate (GHV) as packaged, marketed~~7~~ or
10 manufactured for human consumption, with the exception of legitimate
11 food additive and manufacturing purposes;

12 8. Gamma Valerolactone (GVL) as packaged, marketed~~7~~ or
13 manufactured for human consumption, with the exception of legitimate
14 food additive and manufacturing purposes;

15 9. 1,4 Butanediol (1,4 BD or BDO) as packaged, marketed,
16 manufactured~~7~~ or promoted for human consumption with the exception
17 of legitimate manufacturing purposes; or

18 10. N-ethylpentylone.

19 E. 1. The following industrial uses of Gamma-Butyrolactone,
20 Gamma Hydroxyvalerate, Gamma Valerolactone~~7~~ or 1,4 Butanediol are
21 excluded from all schedules of controlled substances under this
22 title:

23 a. pesticides,

24 b. photochemical etching,

- c. electrolytes of small batteries or capacitors,
- d. viscosity modifiers in polyurethane,
- e. surface etching of metal coated plastics,
- f. organic paint disbursements for water soluble inks,
- g. pH regulators in the dyeing of wool and polyamide fibers,
- h. foundry chemistry as a catalyst during curing,
- 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 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,

1 marketed, manufactured and distributed for legitimate industrial use
2 in a manner that reduces or eliminates the likelihood of abuse.

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

- 6 a. the history and current pattern of abuse,
- 7 b. the name and labeling of the product,
- 8 c. the intended manner of distribution, advertising and
9 promotion of the product, and
- 10 d. other factors as may be relevant to and consistent
11 with the public health and safety.

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

14 F. Any material, compound, mixture~~7~~ or preparation, whether
15 produced directly or indirectly from a substance of vegetable origin
16 or independently by means of chemical synthesis, or by a combination
17 of extraction and chemical synthesis, that contains any quantity of
18 the following substances, or that contains any of their salts,
19 isomers~~7~~ and salts of isomers when the existence of these salts,
20 isomers~~7~~ and salts of isomers is possible within the specific
21 chemical designation:

- 22 1. JWH-004;
- 23 2. JWH-007;
- 24 3. JWH-009;

1	4. JWH-015;
2	5. JWH-016;
3	6. JWH-018;
4	7. JWH-019;
5	8. JWH-020;
6	9. JWH-030;
7	10. JWH-046;
8	11. JWH-047;
9	12. JWH-048;
10	13. JWH-049;
11	14. JWH-050;
12	15. JWH-070;
13	16. JWH-071;
14	17. JWH-072;
15	18. JWH-073;
16	19. JWH-076;
17	20. JWH-079;
18	21. JWH-080;
19	22. JWH-081;
20	23. JWH-082;
21	24. JWH-094;
22	25. JWH-096;
23	26. JWH-098;
24	27. JWH-116;

1	28.	JWH-120;
2	29.	JWH-122;
3	30.	JWH-145;
4	31.	JWH-146;
5	32.	JWH-147;
6	33.	JWH-148;
7	34.	JWH-149;
8	35.	JWH-150;
9	36.	JWH-156;
10	37.	JWH-167;
11	38.	JWH-175;
12	39.	JWH-180;
13	40.	JWH-181;
14	41.	JWH-182;
15	42.	JWH-184;
16	43.	JWH-185;
17	44.	JWH-189;
18	45.	JWH-192;
19	46.	JWH-193;
20	47.	JWH-194;
21	48.	JWH-195;
22	49.	JWH-196;
23	50.	JWH-197;
24	51.	JWH-198;

1	52.	JWH-199;
2	53.	JWH-200;
3	54.	JWH-201;
4	55.	JWH-202;
5	56.	JWH-203;
6	57.	JWH-204;
7	58.	JWH-205;
8	59.	JWH-206;
9	60.	JWH-207;
10	61.	JWH-208;
11	62.	JWH-209;
12	63.	JWH-210;
13	64.	JWH-211;
14	65.	JWH-212;
15	66.	JWH-213;
16	67.	JWH-234;
17	68.	JWH-235;
18	69.	JWH-236;
19	70.	JWH-237;
20	71.	JWH-239;
21	72.	JWH-240;
22	73.	JWH-241;
23	74.	JWH-242;
24	75.	JWH-243;

1	76.	JWH-244;
2	77.	JWH-245;
3	78.	JWH-246;
4	79.	JWH-248;
5	80.	JWH-249;
6	81.	JWH-250;
7	82.	JWH-251;
8	83.	JWH-252;
9	84.	JWH-253;
10	85.	JWH-262;
11	86.	JWH-292;
12	87.	JWH-293;
13	88.	JWH-302;
14	89.	JWH-303;
15	90.	JWH-304;
16	91.	JWH-305;
17	92.	JWH-306;
18	93.	JWH-307;
19	94.	JWH-308;
20	95.	JWH-311;
21	96.	JWH-312;
22	97.	JWH-313;
23	98.	JWH-314;
24	99.	JWH-315;

1	100.	JWH-316;
2	101.	JWH-346;
3	102.	JWH-348;
4	103.	JWH-363;
5	104.	JWH-364;
6	105.	JWH-365;
7	106.	JWH-367;
8	107.	JWH-368;
9	108.	JWH-369;
10	109.	JWH-370;
11	110.	JWH-371;
12	111.	JWH-373;
13	112.	JWH-386;
14	113.	JWH-387;
15	114.	JWH-392;
16	115.	JWH-394;
17	116.	JWH-395;
18	117.	JWH-397;
19	118.	JWH-398;
20	119.	JWH-399;
21	120.	JWH-400;
22	121.	JWH-412;
23	122.	JWH-413;
24	123.	JWH-414;

1	124.	JWH-415;
2	125.	CP-55, 940;
3	126.	CP-47, 497;
4	127.	HU-210;
5	128.	HU-211;
6	129.	WIN-55, 212-2;
7	130.	AM-2201;
8	131.	AM-2233;
9	132.	JWH-018 adamantyl-carboxamide;
10	133.	AKB48;
11	134.	JWH-122 N-(4-pentenyl) analog;
12	135.	MAM2201;
13	136.	URB597;
14	137.	URB602;
15	138.	URB754;
16	139.	UR144;
17	140.	XLR11;
18	141.	A-796,260;
19	142.	STS-135;
20	143.	AB-FUBINACA;
21	144.	AB-PINACA;
22	145.	PB-22;
23	146.	AKB48 N-5-Fluoropentyl;
24	147.	AM1248;

1 148. FUB-PB-22;

2 149. ADB-FUBINACA;

3 150. BB-22;

4 151. 5-Fluoro PB-22; or

5 152. 5-Fluoro AKB-48.

6 G. In addition to those substances listed in subsection F of
7 this section, unless specifically excepted or unless listed in
8 another schedule, any material, compound, mixture, or preparation
9 which contains any quantity of a synthetic cannabinoid found to be
10 in any of the following chemical groups:

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

22 a. 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWH-
23 200),

24 b. 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201),

- c. 1-pentyl-3-(1-naphthoyl)indole (JWH-018),
- d. 1-butyl-3-(1-naphthoyl)indole (JWH-073),
- e. 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (JWH-081),
- f. 1-propyl-2-methyl-3-(1-naphthoyl)indole (JWH-015),
- 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. Naphthylmethylindoles: 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. Naphthylmethylindoles include, but are not limited to, (1-pentylindol-3-yl)(1-naphthyl)methane (JWH-175);

3. Naphthoylpyrroles: any compound containing a 3-(1-naphthoyl)pyrrole structure with or without substitution at the nitrogen atom of the pyrrole 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 pyrrole ring to any extent, and whether or not substituted on the naphthyl group to any extent. Naphthoylpyrroles include, but are not limited to:

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

4. Naphthylideneindenes: any compound containing a 1-(1-naphthylmethylene)indene structure with or without substitution at the 3-position of the indene ring by an alkyl, haloalkyl,

1 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
2 halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
3 morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
4 morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl,
5 phenyl, or halophenyl group, whether or not further substituted on
6 the indene group to any extent, and whether or not substituted on
7 the naphthyl group to any extent. Naphthylmethylindenes include,
8 but are not limited to, (1-[(3-pentyl)-1H-inden-1-
9 ylidene)methyl]naphthalene (JWH-176);

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

- 21 a. 1-pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250),
- 22 b. 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole
23 (RCS-8),
- 24 c. 1-pentyl-3-(2-chlorophenylacetyl)indole (JWH-203),

d. 1-pentyl-3-(2-methylphenylacetyl)indole (JWH-251),

e. 1-pentyl-3-(4-methoxyphenylacetyl)indole (JWH-201), or

f. 1-pentyl-3-(3-methoxyphenylacetyl)indole (JWH-302);

6. Cyclohexylphenols: any compound containing a 2-(3-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~~7~~ 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:

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

12. Carbazole Ketone: Any compound containing (9H-carbazole-3-yl) methanone structure with or without substitution at the nitrogen atom of the carbazole 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~~7~~ or halophenyl group, with substitution at the carbon of the methanone 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 at the carbazole, adamantyl, naphthyl, phenyl, pyrrole, quinolinyl~~7~~ or cycloalkyl rings to any extent. Carbazole Ketones include, but are not limited to, naphthalen-1-yl(9-pentyl-9H-carbazol-3-yl)methanone (EG-018);

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

- 17 a. naphthalen-1-yl(1-pentyl-1H-benzo[d]imidazol-2-
18 1)methanone (JWH-018 benzimidazole analog), or
- 19 b. (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-
20 yl)(naphthalen-1-yl)methanone (FUBIMINA); and

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

SECTION 2. AMENDATORY 63 O.S. 2011, Section 2-210, as last amended by Section 4, Chapter 390, O.S.L. 2017 (63 O.S. Supp. 2019, Section 2-210), is amended to read as follows:

Section 2-210. A. Any material, compound, mixture, or preparation which contains any quantity of the following substances having a potential for abuse associated with a stimulant or depressant effect on the central nervous system:

1. Chloral betaine;
2. Chloral hydrate;
3. Ethchlorvynol;
4. Ethinamate;
5. Meprobamate;
6. Paraldehyde;
7. Petrichloral;
8. Diethylpropion;
9. Phentermine;
10. Pemoline;
11. Chlordiazepoxide;
12. Chlordiazepoxide and its salts, but not including chlordiazepoxide hydrochloride and clidinium bromide or chlordiazepoxide and water-soluble esterified estrogens;
13. Diazepam;
14. Oxazepam;
15. Clorazepate;

- 1 16. Flurazepam and its salts;
- 2 17. Clonazepam;
- 3 18. Barbitol;
- 4 19. Mebutamate;
- 5 20. Methohexital;
- 6 21. Methylphenobarbital;
- 7 22. Phenobarbital;
- 8 23. Fenfluramine;
- 9 24. Pentazocine;
- 10 25. Propoxyphene;
- 11 26. Butorphanol;
- 12 27. Alprazolam;
- 13 28. Halazepam;
- 14 29. Lorazepam;
- 15 30. Prazepam;
- 16 31. Temazepam;
- 17 32. Triazolam;
- 18 33. Carisoprodol;
- 19 34. Dichloralphenazone;
- 20 35. Estazolam;
- 21 36. Eszopiclone;
- 22 37. Midazolam;
- 23 38. Modafinil;
- 24 39. Zaleplon;

40. Zolpidem;
41. Tramadol;
42. Bromazepam;
43. Suvorexant;
44. Phenazepam;
45. Etizolam; ~~or~~
46. Clonazolam; or
47. Gabapentin.

B. 1. The following nonnarcotic substances, which may, under the Federal Food, Drug, and Cosmetic Act (21 U.S.C., Section 301), be lawfully sold over the counter without a prescription, are excluded from all schedules of controlled substances under this title:

- a. Breathe-Aid,
- b. BronCare,
- c. Bronchial Congestion,
- d. Bronkaid Tablets,
- e. Bronkaid Dual Action Caplets,
- f. Bronkotabs,
- g. Bronkolixir,
- h. NeoRespin,
- i. Pazo Hemorrhoid Ointment and Suppositories,
- j. Primatene Tablets,
- k. Primatene "Dual Action" Formula,

1. Quelidrine,
- m. Resp, and
- n. Vatronal Nose Drops.

2. At the request of any person, the Director may exempt any other drug product containing ephedrine from being included as a Schedule IV controlled substance if such product:

- a. is labeled and marketed in a manner consistent with the pertinent OTC tentative final or final monograph issued by the FDA, and
- b. is manufactured and distributed for legitimate medicinal use and in a manner that reduces or eliminates the likelihood of abuse.

3. In making a determination regarding a drug 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 Administrative Procedures Act.

1 5. A list of current drug products meeting exemption
2 requirements under this subsection may be obtained from the Bureau
3 upon written request.

4 C. The Board of Pharmacy may except by rule any compound,
5 mixture, or preparation containing any depressant substance listed
6 in subsection A of this section from the application of all or any
7 part of the Uniform Controlled Dangerous Substances Act, Section 2-
8 101 et seq. of this title, if the compound, mixture, or preparation
9 contains one or more active medicinal ingredients not having a
10 depressant effect on the central nervous system, and if the
11 admixtures are included therein in combinations, quantity,
12 proportion, or concentration that vitiate the potential for abuse of
13 the substances which have a depressant effect on the central nervous
14 system.

15 SECTION 3. This act shall become effective November 1, 2020.

16 COMMITTEE REPORT BY: COMMITTEE ON PUBLIC SAFETY
17 February 10, 2020 - DO PASS AS AMENDED
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