1 ENGROSSED SENATE BILL NO. 1152 By: Standridge of the Senate 2 and 3 Worthen of the House 4 5 An Act relating to the Uniform Controlled Dangerous 6 Substances Act; amending 63 O.S. 2021, Section 2-204, which relates to Schedule I; modifying inclusions; 7 and providing an effective date. 8 9 BE IT ENACTED BY THE PEOPLE OF THE STATE OF OKLAHOMA: 10 63 O.S. 2021, Section 2-204, is 11 SECTION 1. AMENDATORY 12 amended to read as follows: 13 Section 2-204. The controlled substances listed in this section are included in Schedule I and include any material, compound, 14 mixture or preparation that contains any quantity of the following 15 hallucinogenic substances, their salts, isomers and salts of 16 isomers, unless specifically excepted, when the existence of these 17 salts, isomers and salts of isomers is possible within the specific 18 chemical designation. 19 A. Any of the following opiates, including their isomers, 20 esters, ethers, salts, and salts of isomers, esters, and ethers, 21 unless specifically excepted, when the existence of these isomers, 22 esters, ethers, and salts is possible within the specific chemical 23

designation:

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

25. 1 Hydroxypethidine; 26. Ketobemidone; 2 27. Levomoramide; 3 4 28. Levophenacylmorphan; 5 29. Metonitazene; 29. 30. Morpheridine; 6 7 30. 31. Noracymethadol; 31. 32. Norlevorphanol; 8 9 32. 33. Normethadone; 33. 34. Norpipanone; 10 34. 35. Phenadoxone; 11 35. 36. Phenampromide; 12 36. 37. Phenomorphan; 13 37. 38. Phenoperidine; 14 38. 39. Piritramide; 15 39. 40. Proheptazine; 16 40. 41. Properidine; 17 41. 42. Racemoramide; or 18 42. 43. Trimeperidine. 19 B. Any of the following opium derivatives, their salts, 20 isomers, and salts of isomers, unless specifically excepted, when 21 the existence of these salts, isomers, and salts of isomers is 22 possible within the specific chemical designation: 23

24

1. Acetorphine;

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

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

10.

24

Ibogaine;

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

31.

24

B-ketoamphetamine;

```
1
        32.
              2,5-dimethoxy-4-nitroamphetamine;
 2
        33.
              2,5-dimethoxy-4-bromophenethylamine;
 3
        34.
              2,5-dimethoxy-4-chlorophenethylamine;
              2,5-dimethoxy-4-iodoamphetamine;
        35.
 4
 5
        36.
              2,5-dimethoxy-4-iodophenethylamine;
 6
        37.
             2,5-dimethoxy-4-methylphenethylamine;
 7
        38.
              2,5-dimethoxy-4-ethylphenethylamine;
        39.
              2,5-dimethoxy-4-fluorophenethylamine;
 8
 9
        40.
             2,5-dimethoxy-4-nitrophenethylamine;
              2,5-dimethoxy-4-ethylthio-phenethylamine;
10
        41.
        42.
             2,5-dimethoxy-4-isopropylthio-phenethylamine;
11
              2,5-dimethoxy-4-propylthio-phenethylamine;
12
        43.
13
        44.
              2,5-dimethoxy-4-cyclopropylmethylthio-phenethylamine;
        45.
              2,5-dimethoxy-4-tert-butylthio-phenethylamine;
14
        46.
              2,5-dimethoxy-4-(2-fluoroethylthio)-phenethylamine;
15
        47.
              5-methoxy-N, N-dimethyltryptamine;
16
        48.
             N-methyltryptamine;
17
             A-ethyltryptamine;
        49.
18
             A-methyltryptamine;
        50.
19
             N, N-diethyltryptamine;
        51.
20
        52.
             N, N-diisopropyltryptamine;
21
             N, N-dipropyltryptamine;
        53.
22
             5-methoxy-a-methyltryptamine;
        54.
23
              4-hydroxy-N, N-diethyltryptamine;
        55.
24
```

```
56.
 1
              4-hydroxy-N, N-diisopropyltryptamine;
 2
        57.
              5-methoxy-N, N-diisopropyltryptamine;
        58.
              4-hydroxy-N-isopropyl-N-methyltryptamine;
 3
        59.
              3,4-Methylenedioxymethcathinone (Methylone);
 4
              3,4-Methylenedioxypyrovalerone (MDPV);
 5
        60.
        61.
              4-Methylmethcathinone (Mephedrone);
 6
 7
        62.
             4-methoxymethcathinone;
        63.
             4-Fluoromethcathinone;
 8
        64.
 9
             3-Fluoromethcathinone;
        65.
             1-(8-bromobenzo 1,2-b;4,5-b' difuran-4-yl)-2-aminopropane;
10
        66.
             2,5-Dimethoxy-4-chloroamphetamine;
11
12
        67.
             4-Methylethcathinone;
13
        68.
             Pyrovalerone;
        69.
             N, N-diallyl-5-methoxytryptamine;
14
        70.
              3,4-Methylenedioxy-N-ethylcathinone (Ethylone);
15
        71.
             B-keto-N-Methylbenzodioxolylbutanamine (Butylone);
16
        72.
             B-keto-Methylbenzodioxolylpentanamine (Pentylone);
17
        73.
             Alpha-Pyrrolidinopentiophenone;
18
        74.
             4-Fluoroamphetamine;
19
        75.
             Pentedrone;
20
        76.
             4'-Methyl-a-pyrrolidinohexaphenone;
21
        77.
             2,5-dimethoxy-4-(n)-propylphenethylamine;
22
        78.
             2,5-dimethoxyphenethylamine;
23
        79.
             1,4-Dibenzylpiperazine;
24
```

```
1
        80.
             N, N-Dimethylamphetamine;
 2
             4-Fluoromethamphetamine;
        81.
             4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine
 3
        82.
    (25C-NBOMe);
 4
 5
             4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine
    (25I-NBOMe);
 6
 7
             4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine
        84.
    (25B-NBOMe);
 8
 9
        85.
            1-(4-Fluorophenyl)piperazine;
        86.
            Methoxetamine;
10
             3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N-
11
        87.
12
    methylbenzamide;
13
        88.
             N-ethyl hexadrone;
        89.
             Isopropyl-U-47700;
14
        90.
             Para-fluorobutyrl fentanyl;
15
             Fluoro isobutryrl fentanyl;
        91.
16
        92.
             3-Hydroxy Phencyclidine (PCP); or
17
             3-methoxy Phencyclidine (PCP);
        93.
18
        94. Flualprazolam; or
19
        95. Flubromazolam.
20
            Unless specifically excepted or unless listed in a different
21
    schedule, any material, compound, mixture, or preparation which
22
    contains any quantity of the following substances having stimulant
23
    or depressant effect on the central nervous system:
24
```

- 1 1. Fenethylline;
- 2 2. Mecloqualone;
- 3. N-ethylamphetamine;
 - 4. Methaqualone;

11

12

13

14

15

16

17

18

19

- 5. Gamma-Hydroxybutyric Acid, also known as GHB, gamma6 hydroxybutyrate, 4-hydroxybutyrate, 4-hydroxybutanoic acid, sodium
 7 oxybate, and sodium oxybutyrate;
- 8 6. Gamma-Butyrolactone (GBL) as packaged, marketed,
 9 manufactured or promoted for human consumption, with the exception
 10 of legitimate food additive and manufacturing purposes;
 - 7. Gamma Hydroxyvalerate (GHV) as packaged, marketed, or manufactured for human consumption, with the exception of legitimate food additive and manufacturing purposes;
 - 8. Gamma Valerolactone (GVL) as packaged, marketed, or manufactured for human consumption, with the exception of legitimate food additive and manufacturing purposes;
 - 9. 1,4 Butanediol (1,4 BD or BDO) as packaged, marketed, manufactured, or promoted for human consumption with the exception of legitimate manufacturing purposes; or
 - 10. N-ethylpentylone.
- E. 1. The following industrial uses of Gamma-Butyrolactone,
 Gamma Hydroxyvalerate, Gamma Valerolactone, or 1,4 Butanediol are
 excluded from all schedules of controlled substances under this
 title:

1	a.	pesticides,
2	b.	photochemical etching,
3	С.	electrolytes of small batteries or capacitors,
4	d.	viscosity modifiers in polyurethane,
5	е.	surface etching of metal coated plastics,
6	f.	organic paint disbursements for water soluble inks,
7	g.	pH regulators in the dyeing of wool and polyamide
8		fibers,
9	h.	foundry chemistry as a catalyst during curing,
10	i.	curing agents in many coating systems based on
11		urethanes and amides,
12	j.	additives and flavoring agents in food, confectionary,
13		and beverage products,
14	k.	synthetic fiber and clothing production,
15	1.	tetrahydrofuran production,
16	m.	gamma butyrolactone production,
17	n.	polybutylene terephthalate resin production,
18	٥.	polyester raw materials for polyurethane elastomers
19		and foams,
20	p.	coating resin raw material, and
21	ď.	as an intermediate in the manufacture of other
22		chemicals and pharmaceuticals.
23	2. At th	e request of any person, the Director may exempt any
24	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. 1 JWH-007; 2 3. JWH-009; 3 4. JWH-015; 4 5. JWH-016; 6. JWH-018; 5 6 7. JWH-019; 7 8. JWH-020; 9. 8 JWH-030; 9 10. JWH-046; 11. JWH-047;10 12. JWH-048; 11 13. JWH-049;12 13 14. JWH-050; 15. JWH-070; 14 16. JWH-071; 15 17. JWH-072; 16 18. JWH-073; 17 19. JWH-076; 18 20. JWH-079; 19 21. JWH-080; 20 22. JWH-081; 21 23. JWH-082; 22 JWH-094; 24. 23 25. JWH-096;

1	26.	JWH-098;
2	27.	JWH-116;
3	28.	JWH-120;
4	29.	JWH-122;
5	30.	JWH-145;
6	31.	JWH-146;
7	32.	JWH-147;
8	33.	JWH-148;
9	34.	JWH-149;
10	35.	JWH-150;
11	36.	JWH-156;
12	37.	JWH-167;
13	38.	JWH-175;
14	39.	JWH-180;
15	40.	JWH-181;
16	41.	JWH-182;
17	42.	JWH-184;
18	43.	JWH-185;
19	44.	JWH-189;
20	45.	JWH-192;
21	46.	JWH-193;
22	47.	JWH-194;
23	48.	JWH-195;
24	49.	JWH-196;

1	50.	JWH-197;
2	51.	JWH-198;
3	52.	JWH-199;
4	53.	JWH-200;
5	54.	JWH-201;
6	55.	JWH-202;
7	56.	JWH-203;
8	57.	JWH-204;
9	58.	JWH-205;
10	59.	JWH-206;
11	60.	JWH-207;
12	61.	JWH-208;
13	62.	JWH-209;
14	63.	JWH-210;
15	64.	JWH-211;
16	65.	JWH-212;
17	66.	JWH-213;
18	67.	JWH-234;
19	68.	JWH-235;
20	69.	JWH-236;
21	70.	JWH-237;
22	71.	JWH-239;
23	72.	JWH-240;
24	73.	JWH-241;

1	74.	JWH-242;
2	75.	JWH-243;
3	76.	JWH-244;
4	77.	JWH-245;
5	78.	JWH-246;
6	79.	JWH-248;
7	80.	JWH-249;
8	81.	JWH-250;
9	82.	JWH-251;
10	83.	JWH-252;
11	84.	JWH-253;
12	85.	JWH-262;
13	86.	JWH-292;
14	87.	JWH-293;
15	88.	JWH-302;
16	89.	JWH-303;
17	90.	JWH-304;
18	91.	JWH-305;
19	92.	JWH-306;
20	93.	JWH-307;
21	94.	JWH-308;
22	95.	JWH-311;
23	96.	JWH-312;
24	97.	JWH-313;

1	98.	JWH-314;
2	99.	JWH-315;
3	100.	JWH-316;
4	101.	JWH-346;
5	102.	JWH-348;
6	103.	JWH-363;
7	104.	JWH-364;
8	105.	JWH-365;
9	106.	JWH-367;
10	107.	JWH-368;
11	108.	JWH-369;
12	109.	JWH-370;
13	110.	JWH-371;
14	111.	JWH-373;
15	112.	JWH-386;
16	113.	JWH-387;
17	114.	JWH-392;
18	115.	JWH-394;
19	116.	JWH-395;
20	117.	JWH-397;
21	118.	JWH-398;
22	119.	JWH-399;
23	120.	JWH-400;
24	121.	JWH-412;

```
1
        122. JWH-413;
 2
        123. JWH-414;
 3
        124. JWH-415;
 4
        125. CP-55, 940;
        126. CP-47, 497;
 5
 6
        127. HU-210;
 7
        128. HU-211;
        129. WIN-55, 212-2;
 8
 9
        130. AM-2201;
        131. AM-2233;
10
              JWH-018 adamantyl-carboxamide;
        132.
11
        133. AKB48;
12
13
        134.
              JWH-122 N-(4-pentenyl) analog;
        135. MAM2201;
14
        136. URB597;
15
        137. URB602;
16
        138. URB754;
17
        139. UR144;
18
        140. XLR11;
19
        141. A-796,260;
20
        142. STS-135;
21
        143. AB-FUBINACA;
22
        144. AB-PINACA;
23
        145. PB-22;
24
```

```
1
        146.
              AKB48 N-5-Fluorpentyl;
 2
        147.
              AM1248;
        148.
              FUB-PB-22;
 3
        149.
              ADB-FUBINACA;
 4
 5
        150.
              BB-22;
        151.
              5-Fluoro PB-22; or
 6
 7
        152.
              5-Fluoro AKB-48.
            In addition to those substances listed in subsection F of
 8
 9
    this section, unless specifically excepted or unless listed in
    another schedule, any material, compound, mixture, or preparation
10
    which contains any quantity of a synthetic cannabinoid found to be
11
    in any of the following chemical groups:
12
13
        1. Naphthoylindoles: any compound containing a 3-(1-
    naphthoyl)indole structure with or without substitution at the
14
    nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,
15
    alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-
16
17
    (N-methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-
    2-pyrrolidinyl) methyl, 1-(N-methyl-3- morpholinyl) methyl,
18
```

19 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 20 halophenyl group, whether or not further substituted on the indole

ring to any extent, and whether or not substituted on the naphthyl

22 ring to any extent. Naphthoylindoles include, but are not limited

23 to:

24

```
1
                   1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWH-
             a.
 2
                   200).
                   1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201),
 3
             b.
                   1-pentyl-3-(1-naphthoyl)indole (JWH-018),
 4
             C.
 5
             d.
                   1-butyl-3-(1-naphthoyl)indole (JWH-073),
                   1-pentyl-3-(4-methoxy-1-naphthoyl)indole (JWH-081),
 6
             е.
 7
             f.
                   1-propyl-2-methyl-3-(1-naphthoyl)indole (JWH-015),
                   1-hexyl-3-(1-naphthoyl)indole (JWH-019),
 8
             g.
 9
             h.
                   1-pentyl-3-(4-methyl-1-naphthoyl)indole (JWH-122),
10
             i.
                   1-pentyl-3-(4-ethyl-1-naphthoyl)indole (JWH-210),
                   1-pentyl-3-(4-chloro-1-naphthoyl)indole (JWH-398),
11
             j.
                   1-pentyl-2-methyl-3-(1-naphthoyl)indole (JWH-007),
12
             k.
             1.
                   1-pentyl-3-(7-methoxy-1-naphthoyl)indole (JWH-164),
13
                   1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole
14
             m.
                   (JWH-098),
15
                   1-pentyl-3-(4-fluoro-1-naphthoyl)indole (JWH-412),
16
             n.
                   1-[1-(N-methyl-2-piperidinyl) methyl]-3-(1-
17
             Ο.
                   naphthoyl) indole (AM-1220),
18
                   1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole
19
             р.
                   (MAM-2201), or
20
                   1-(4-cyanobutyl)-3-(1-naphthoyl) indole (AM-2232);
21
             q.
            Naphthylmethylindoles: any compound containing a 1H-indol-3-
22
    yl-(1-naphthyl) methane structure with or without substitution at the
23
24
    nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,
```

```
1
    alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-
 2
    (N-methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-
 3
    2-pyrrolidinyl) methyl, 1-(N-methyl-3- morpholinyl) methyl,
 4
    (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
 5
    halophenyl group, whether or not further substituted on the indole
    ring to any extent, and whether or not substituted on the naphthyl
 6
    ring to any extent. Naphthylmethylindoles include, but are not
 7
    limited to, (1-pentylindol-3-yl)(1-naphthyl)methane (JWH-175);
 8
 9
        3. Naphthoylpyrroles: any compound containing a 3-(1-
10
    naphthoyl)pyrrole structure with or without substitution at the
11
    nitrogen atom of the pyrrole ring by an alkyl, haloalkyl,
12
    cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
13
    halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
    morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
14
    morpholinyl) methyl, (tetrahydropyran-4-yl) methyl, 1-methylazepanyl,
15
    phenyl, or halophenyl group, whether or not further substituted on
16
17
    the pyrrole ring to any extent, and whether or not substituted on
    the naphthyl group to any extent. Naphthoylpyrroles include, but
18
    are not limited to:
19
                  1-hexyl-2-phenyl-4-(1-naphthoyl)pyrrole (JWH-147),
20
             a.
             b.
                  1-pentyl-5-(2-methylphenyl)-3-(1-naphthoyl)pyrrole
21
                  (JWH - 370),
22
                  1-pentyl-3-(1-naphthoyl)pyrrole (JWH-030), or
23
             C.
```

1-hexyl-5-phenyl-3-(1-naphthoyl)pyrrole (JWH-147);

d.

```
1
            Naphthylideneindenes: any compound containing a 1-(1-
    naphthylmethylene) indene structure with or without substitution at
 2
    the 3-position of the indene ring by an alkyl, haloalkyl,
 3
    cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
 4
 5
    halobenzyl, 1-(N-methyl-2-piperidinyl) methyl, 2-(4-
    morpholinyl) ethyl, 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
 6
    morpholinyl) methyl, (tetrahydropyran-4-yl) methyl, 1-methylazepanyl,
 7
    phenyl, or halophenyl group, whether or not further substituted on
 8
 9
    the indene group to any extent, and whether or not substituted on
10
    the naphthyl group to any extent. Naphthylmethylindenes include,
    but are not limited to, (1-[(3-pentyl)-1H-inden-1-
11
    ylidene)methyl]naphthalene (JWH-176);
12
```

- 5. Phenylacetylindoles: any compound containing a 3-13 phenylacetylindole structure with or without substitution at the 14 nitrogen atom of the indole ring by alkyl, haloalkyl, cyanoalkyl, 15 alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-16 17 (N-methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3- morpholinyl) methyl, 18 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 19 halophenyl group, whether or not further substituted on the indole 20 ring to any extent, and whether or not substituted on the phenyl 21 ring to any extent. Phenylacetylindoles include, but are not 22 limited to: 23
 - a. 1-pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250),

```
1
             b.
                  1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole
 2
                   (RCS-8),
                  1-pentyl-3-(2-chlorophenylacetyl)indole (JWH-203),
 3
             C.
                  1-pentyl-3-(2-methylphenylacetyl)indole (JWH-251),
 4
             d.
 5
                  1-pentyl-3-(4-methoxyphenylacetyl)indole (JWH-201), or
             е.
             f.
                  1-pentyl-3-(3-methoxyphenylacetyl)indole (JWH-302);
 6
 7
            Cyclohexylphenols: any compound containing a 2-(3-
        6.
    hydroxycyclohexyl) phenol structure with or without substitution at
 8
 9
    the 5-position of the phenolic ring by an alkyl, haloalkyl,
    cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
10
    halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
11
    morpholinyl) ethyl, 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
12
13
    morpholinyl) methyl, (tetrahydropyran-4-yl) methyl, 1-methylazepanyl,
    phenyl, or halophenyl group, and whether or not further substituted
14
    on the cyclohexyl ring to any extent. Cyclohexylphenols include,
15
    but are not limited to:
16
                  5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-
17
             a.
                  hydroxycyclohexyl]-phenol (CP-47,497),
18
             b.
                  5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-
19
                  phenol (cannabicyclohexanol; CP-47,497 C8 homologue),
20
                  or
21
                  5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-
22
             C.
                  hydroxypropyl)cyclohexyl]-phenol (CP 55, 940);
23
```

```
1
        7. Benzoylindoles: any compound containing a 3-(benzoyl)indole
    structure with or without substitution at the nitrogen atom of the
 2
    indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
 3
    cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-
 4
 5
    2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-
    pyrrolidinyl) methyl, 1-(N-methyl-3- morpholinyl) methyl,
 6
    (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
 7
    halophenyl group, whether or not further substituted on the indole
 8
 9
    ring to any extent, and whether or not substituted on the phenyl
    group to any extent. Benzoylindoles include, but are not limited
10
11
    to:
                  1-pentyl-3-(4-methoxybenzoyl)indole (RCS-4),
12
             a.
             b.
                  1-[2-(4-morpholinyl)ethyl]-2-methyl-3-(4-
13
                  methoxybenzoyl) indole (Pravadoline or WIN 48, 098),
14
```

- 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-methyl2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl,

16

17

18

19

20

21

22

23

```
1
   (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
   halophenyl group, whether or not further substituted in the indole
2
   ring to any extent and whether or not substituted in the
3
   cyclopropoyl ring to any extent. Cyclopropoylindoles include, but
4
5
   are not limited to:
                 1-pentyl-3-(2,2,3,3-tetramethylcyclopropoyl)indole
6
            a.
                  (UR-144),
7
8
            b.
```

- b. 1-(5-chloropentyl)-3-(2,2,3,3tetramethylcyclopropoyl)indole (5Cl-UR-144), or
- c. 1-(5-fluoropentyl)-3-(2,2,3,3tetramethylcyclopropoyl)indole (XLR11);
- Indole Amides: Any compound containing a 1H-Indole-3-12 9. 13 carboxamide structure with or without substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, 14 cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-15 2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-16 pyrrolidinyl) methyl, 1-(N-methyl-3- morpholinyl) methyl, 17 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 18 halophenyl group, whether or not substituted at the carboxamide 19 group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl, 20 cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-21 1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-22 dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not 23 further substituted in the indole, adamantyl, naphthyl, phenyl, 24

10

```
1
    pyrrole, quninolinyl, or cycloalkyl rings to any extent.
    Amides include, but are not limited to:
 2
                  N-(1-adamantyl)-1-pentyl-1H-indole-3-carboxamide
 3
             a.
                   (2NE1),
 4
 5
             b.
                  N-(1-adamantyl)-1-(5-fluoropentyl-1H-indole-3-
                  carboxamide (STS-135),
 6
                  N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
 7
             C.
                  indole-3-carboxamide (ADBICA),
 8
 9
             d.
                  N-(1-amino-3, 3-dimethyl-1-oxobutan-2-yl)-1-(5-
                  fluoropentyl)-1H-indole-3-carboxamide (5F-ADBICA),
10
                  N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide
11
             е.
12
                  (NNE1),
             f.
                  1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-
13
                  carboxamide (5F-NNE1),
14
                  N-benzyl-1-pentyl-1H-indole-3-carboxamide (SDB-006),
15
             g.
16
                  or
             h.
                  N-benzyl-1-(5-fluoropentyl)-1H-indole-3-carboxamide
17
                   (5F-SDB-006);
18
        10.
             Indole Esters: Any compound containing a 1H-Indole-3-
19
    carboxylate structure with or without substitution at the nitrogen
20
    atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
21
    cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-
22
```

2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-

pyrrolidinyl) methyl, 1-(N-methyl-3-morpholinyl) methyl,

23

```
1
   (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
   halophenyl group, whether or not substituted at the carboxylate
2
   group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,
3
   cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-
4
5
   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
6
   further substituted in the indole, adamantyl, naphthyl, phenyl,
7
   pyrrole, quinolinyl, or cycloalkyl rings to any extent. Indole
8
```

Esters include, but are not limited to:

- a. quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB22),
- b. quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3carboxylate (5F-PB-22),
- c. quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3carboxylate (BB-22),
- d. naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3carboxylate (FDU-PB-22), or
- e. naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3carboxylate (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-

9

10

11

12

13

14

15

16

17

18

19

20

21

22

23

- 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:
 - 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 (AB001);
- Carbazole Ketone: Any compound containing (9H-carbazole-3-11 12. yl) methanone structure with or without substitution at the nitrogen 12 13 atom of the carbazole ring by an alkyl, haloalkyl, cyanoalkyl, 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 17 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, with substitution at the carbon of the methanone 18 group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl, 19 cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-20 1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-21 dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not 22 further substituted at the carbazole, adamantyl, naphthyl, phenyl, 23 pyrrole, quinolinyl, or cycloalkyl rings to any extent. Carbazole 24

8

9

```
Ketones include, but are not limited to, naphthalen-1-yl(9-pentyl-
 1
    9H-carbazol-3-yl)methanone (EG-018);
 2
        13.
             Benzimidazole Ketone: Any compound containing
 3
 4
    (benzimidazole-2-yl) methanone structure with or without
 5
    substitution at either nitrogen atom of the benzimidazole ring by an
    alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
 6
    cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-
 7
    piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-
 8
 9
    pyrrolidinyl) methyl, 1-(N-methyl-3-morpholinyl) methyl,
    (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
10
    halophenyl group, with substitution at the carbon of the methanone
11
    group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,
12
13
    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-
14
    dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not
15
    further substituted in the benzimidazole, adamantyl, naphthyl,
16
17
    phenyl, pyrrole, quinolinyl, or cycloalkyl rings to any extent.
    Benzimidazole Ketones include, but are not limited to:
18
                  naphthalen-1-yl(1-pentyl-1H-benzo[d]imidazol-2-
19
             a.
                  1) methanone (JWH-018 benzimidazole analog), or
20
             b.
                  (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-
21
                  yl) (naphthalen-1-yl) methanone (FUBIMINA); and
22
23
```

- 14. Modified by Replacement: any compound defined in this subsection that is modified by replacement of a carbon with nitrogen in the indole, naphthyl, indene, benzimidazole, or carbazole ring.
- Any prescription drug approved by the federal Food and Drug Administration under the provisions of Section 505 of the Federal Food, Drug and Cosmetic Act, Title 21 of the United States Code, Section 355, that is designated, rescheduled or deleted as a controlled substance under federal law by the United States Drug Enforcement Administration shall be excluded from Schedule I and shall be prescribed, distributed, dispensed or used in accordance with federal law upon the issuance of a notice, final rule or interim final rule by the United States Drug Enforcement Administration designating, rescheduling or deleting as a controlled substance such a drug product under federal law, unless and until the Board of Pharmacy takes action pursuant to Section 2-201 of this title. If the Board of Pharmacy does not take action pursuant to Section 2-201 of this title, the drug product shall be deemed to be designated, rescheduled or deleted as a controlled substance in accordance with federal law and in compliance with the Uniform Controlled Dangerous Substances Act.

SECTION 2. This act shall become effective November 1, 2022.

22

1

2

3

4

5

6

7

9

10

11

12

13

14

15

16

17

18

19

20

21

23

1	Passed the Senate the 28th day of February, 2022.
2	
3	
4	Presiding Officer of the Senate
5	Passed the House of Representatives the day of,
6	2022.
7	
8	
9	Presiding Officer of the House of Representatives
10	
11	
12	
13	
14	
15	
16	
17	
18	
19	
20	
21	
22	
23	
24	