

1 AN ACT concerning criminal law.

2 **Be it enacted by the People of the State of Illinois,**  
3 **represented in the General Assembly:**

4 Section 5. The Illinois Controlled Substances Act is  
5 amended by changing Section 204 as follows:

6 (720 ILCS 570/204) (from Ch. 56 1/2, par. 1204)

7 Sec. 204. (a) The controlled substances listed in this  
8 Section are included in Schedule I.

9 (b) Unless specifically excepted or unless listed in  
10 another schedule, any of the following opiates, including  
11 their isomers, esters, ethers, salts, and salts of isomers,  
12 esters, and ethers, whenever the existence of such isomers,  
13 esters, ethers and salts is possible within the specific  
14 chemical designation:

15 (1) Acetylmethadol;

16 (1.1) Acetyl-alpha-methylfentanyl

17 (N-[1-(1-methyl-2-phenethyl)-

18 4-piperidinyl]-N-phenylacetamide);

19 (2) Allylprodine;

20 (3) Alphacetylmethadol, except

21 levo-alphacetylmethadol (also known as levo-alpha-

22 acetylmethadol, levomethadyl acetate, or LAAM);

23 (4) Alphameprodine;

- 1 (5) Alphamethadol;
- 2 (6) Alpha-methylfentanyl
- 3 (N-(1-alpha-methyl-beta-phenyl) ethyl-4-piperidyl)
- 4 propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-
- 5 propanilido) piperidine;
- 6 (6.1) Alpha-methylthiofentanyl
- 7 (N-[1-methyl-2-(2-thienyl)ethyl-
- 8 4-piperidinyl]-N-phenylpropanamide);
- 9 (7) 1-methyl-4-phenyl-4-propionoxypiperidine (MPPP);
- 10 (7.1) PEPAP
- 11 (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 12 (8) Benzethidine;
- 13 (9) Betacetylmethadol;
- 14 (9.1) Beta-hydroxyfentanyl
- 15 (N-[1-(2-hydroxy-2-phenethyl)-
- 16 4-piperidinyl]-N-phenylpropanamide);
- 17 (10) Betameprodine;
- 18 (11) Betamethadol;
- 19 (12) Betaprodine;
- 20 (13) Clonitazene;
- 21 (14) Dextromoramide;
- 22 (15) Diampromide;
- 23 (16) Diethylthiambutene;
- 24 (17) Difenoazin;
- 25 (18) Dimenoxadol;
- 26 (19) Dimepheptanol;

- 1 (20) Dimethylthiambutene;
- 2 (21) Dioxaphetylbutyrate;
- 3 (22) Dipipanone;
- 4 (23) Ethylmethylthiambutene;
- 5 (24) Etonitazene;
- 6 (25) Etoxeridine;
- 7 (26) Furethidine;
- 8 (27) Hydroxypethidine;
- 9 (28) Ketobemidone;
- 10 (29) Levomoramide;
- 11 (30) Levophenacymorphan;
- 12 (31) 3-Methylfentanyl
- 13 (N-[3-methyl-1-(2-phenylethyl)-
- 14 4-piperidyl]-N-phenylpropanamide);
- 15 (31.1) 3-Methylthiofentanyl
- 16 (N-[(3-methyl-1-(2-thienyl)ethyl-
- 17 4-piperidinyl]-N-phenylpropanamide);
- 18 (32) Morpheridine;
- 19 (33) Noracymethadol;
- 20 (34) Norlevorphanol;
- 21 (35) Normethadone;
- 22 (36) Norpipanone;
- 23 (36.1) Para-fluorofentanyl
- 24 (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-
- 25 4-piperidinyl]propanamide);
- 26 (37) Phenadoxone;

- 1 (38) Phenampromide;
- 2 (39) Phenomorphan;
- 3 (40) Phenoperidine;
- 4 (41) Piritramide;
- 5 (42) Proheptazine;
- 6 (43) Properidine;
- 7 (44) Propiram;
- 8 (45) Racemoramide;
- 9 (45.1) Thiofentanyl
- 10 (N-phenyl-N-[1-(2-thienyl)ethyl-
- 11 4-piperidinyl]-propanamide);
- 12 (46) Tilidine;
- 13 (47) Trimeperidine;
- 14 (48) Beta-hydroxy-3-methylfentanyl (other name:
- 15 N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-
- 16 N-phenylpropanamide);
- 17 (49) Furanyl fentanyl (FU-F);
- 18 (50) Butyryl fentanyl;
- 19 (51) Valeryl fentanyl;
- 20 (52) Acetyl fentanyl;
- 21 (53) Beta-hydroxy-thiofentanyl;
- 22 (54) 3,4-dichloro-N-[2-
- 23 (dimethylamino)cyclohexyl]-N-
- 24 methylbenzamide (U-47700);
- 25 (55) 4-chloro-N-[1-[2-
- 26 (4-nitrophenyl)ethyl]-2-piperidinylidene]-

1 benzenesulfonamide (W-18);  
2 (56) 4-chloro-N-[1-(2-phenylethyl)  
3 -2-piperidinylidene]-benzenesulfonamide (W-15);  
4 (57) acrylfentanyl (acryloylfentanyl).

5 (c) Unless specifically excepted or unless listed in  
6 another schedule, any of the following opium derivatives, its  
7 salts, isomers and salts of isomers, whenever the existence of  
8 such salts, isomers and salts of isomers is possible within  
9 the specific chemical designation:

- 10 (1) Acetorphine;
- 11 (2) Acetyldihydrocodeine;
- 12 (3) Benzylmorphine;
- 13 (4) Codeine methylbromide;
- 14 (5) Codeine-N-Oxide;
- 15 (6) Cyprenorphine;
- 16 (7) Desomorphine;
- 17 (8) Diacetyldihydromorphine (Dihydroheroin);
- 18 (9) Dihydromorphine;
- 19 (10) Drotebanol;
- 20 (11) Etorphine (except hydrochloride salt);
- 21 (12) Heroin;
- 22 (13) Hydromorphenol;
- 23 (14) Methyldesorphine;
- 24 (15) Methyldihydromorphine;
- 25 (16) Morphine methylbromide;
- 26 (17) Morphine methylsulfonate;

- 1 (18) Morphine-N-Oxide;
- 2 (19) Myrophine;
- 3 (20) Nicocodeine;
- 4 (21) Nicomorphine;
- 5 (22) Normorphine;
- 6 (23) Pholcodine;
- 7 (24) Thebacon.

8 (d) Unless specifically excepted or unless listed in  
9 another schedule, any material, compound, mixture, or  
10 preparation which contains any quantity of the following  
11 hallucinogenic substances, or which contains any of its salts,  
12 isomers and salts of isomers, whenever the existence of such  
13 salts, isomers, and salts of isomers is possible within the  
14 specific chemical designation (for the purposes of this  
15 paragraph only, the term "isomer" includes the optical,  
16 position and geometric isomers):

- 17 (1) 3,4-methylenedioxyamphetamine  
18 (alpha-methyl,3,4-methylenedioxyphenethylamine,  
19 methylenedioxyamphetamine, MDA);  
20 (1.1) Alpha-ethyltryptamine  
21 (some trade or other names: etryptamine;  
22 MONASE; alpha-ethyl-1H-indole-3-ethanamine;  
23 3-(2-aminobutyl)indole; a-ET; and AET);  
24 (2) 3,4-methylenedioxymethamphetamine (MDMA);  
25 (2.1) 3,4-methylenedioxy-N-ethylamphetamine  
26 (also known as: N-ethyl-alpha-methyl-

1 3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE,  
2 and MDEA);

3 (2.2) N-Benzylpiperazine (BZP);

4 (2.2-1) Trifluoromethylphenylpiperazine (TFMPP);

5 (3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);

6 (4) 3,4,5-trimethoxyamphetamine (TMA);

7 (5) (Blank);

8 (6) Diethyltryptamine (DET);

9 (7) Dimethyltryptamine (DMT);

10 (7.1) 5-Methoxy-diallyltryptamine;

11 (8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);

12 (9) Ibogaine (some trade and other names:  
13 7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-  
14 6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b]  
15 indole; Tabernanthe iboga);

16 (10) Lysergic acid diethylamide;

17 (10.1) Salvinorin A;

18 (10.5) Salvia divinorum (meaning all parts of the  
19 plant presently classified botanically as Salvia  
20 divinorum, whether growing or not, the seeds thereof, any  
21 extract from any part of that plant, and every compound,  
22 manufacture, salts, isomers, and salts of isomers whenever  
23 the existence of such salts, isomers, and salts of isomers  
24 is possible within the specific chemical designation,  
25 derivative, mixture, or preparation of that plant, its  
26 seeds or extracts);

- 1 (11) 3,4,5-trimethoxyphenethylamine (Mescaline);
- 2 (12) Peyote (meaning all parts of the plant presently  
3 classified botanically as *Lophophora williamsii* Lemaire,  
4 whether growing or not, the seeds thereof, any extract  
5 from any part of that plant, and every compound,  
6 manufacture, salts, derivative, mixture, or preparation of  
7 that plant, its seeds or extracts);
- 8 (13) N-ethyl-3-piperidyl benzilate (JB 318);
- 9 (14) N-methyl-3-piperidyl benzilate;
- 10 (14.1) N-hydroxy-3,4-methylenedioxyamphetamine  
11 (also known as N-hydroxy-alpha-methyl-  
12 3,4(methylenedioxy)phenethylamine and N-hydroxy MDA);
- 13 (15) Parahexyl; some trade or other names:  
14 3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-  
15 dibenzo (b,d) pyran; Synhexyl;
- 16 (16) Psilocybin;
- 17 (17) Psilocyn;
- 18 (18) Alpha-methyltryptamine (AMT);
- 19 (19) 2,5-dimethoxyamphetamine  
20 (2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA);
- 21 (20) 4-bromo-2,5-dimethoxyamphetamine  
22 (4-bromo-2,5-dimethoxy-alpha-methylphenethylamine;  
23 4-bromo-2,5-DMA);
- 24 (20.1) 4-Bromo-2,5 dimethoxyphenethylamine.  
25 Some trade or other names: 2-(4-bromo-  
26 2,5-dimethoxyphenyl)-1-aminoethane;



1 alpha-desmethyl DOB, 2CB, Nexus;

2 (21) 4-methoxyamphetamine

3 (4-methoxy-alpha-methylphenethylamine;

4 paramethoxyamphetamine; PMA);

5 (22) (Blank);

6 (23) Ethylamine analog of phencyclidine.

7 Some trade or other names:

8 N-ethyl-1-phenylcyclohexylamine,

9 (1-phenylcyclohexyl) ethylamine,

10 N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE;

11 (24) Pyrrolidine analog of phencyclidine. Some trade

12 or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy,

13 PHP;

14 (25) 5-methoxy-3,4-methylenedioxy-amphetamine;

15 (26) 2,5-dimethoxy-4-ethylamphetamine

16 (another name: DOET);

17 (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine

18 (another name: TCPy);

19 (28) (Blank);

20 (29) Thiophene analog of phencyclidine (some trade

21 or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine;

22 2-thienyl analog of phencyclidine; TPCP; TCP);

23 (29.1) Benzothiophene analog of phencyclidine. Some

24 trade or other names: BTCP or benocyclidine;

25 (29.2) 3-Methoxyphencyclidine (3-MeO-PCP);

26 (30) Bufotenine (some trade or other names:

1 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole;  
2 3-(2-dimethylaminoethyl)-5-indolol;  
3 5-hydroxy-N,N-dimethyltryptamine;  
4 N,N-dimethylserotonin; mappine);  
5 (31) (Blank);  
6 (32) (Blank);  
7 (33) (Blank);  
8 (34) (Blank);  
9 (34.5) (Blank);  
10 (35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-  
11 (2-methyloctan-2-yl)-6a,7,  
12 10,10a-tetrahydrobenzo[c]chromen-1-ol  
13 Some trade or other names: HU-210;  
14 (35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-  
15 dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-  
16 tetrahydrobenzo[c]chromen-1-ol, its isomers,  
17 salts, and salts of isomers; Some trade or other  
18 names: HU-210, Dexanabinol;  
19 (36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-  
20 6,6-dimethyl-3-(2-methyloctan-2-yl)-  
21 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol  
22 Some trade or other names: HU-211;  
23 (37) (Blank);  
24 (38) (Blank);  
25 (39) (Blank);  
26 (40) (Blank);

1 (41) (Blank);

2 (42) Any compound structurally derived from  
3 3-(1-naphthoyl)indole or  
4 1H-indol-3-yl-(1-naphthyl)methane by substitution at the  
5 nitrogen atom of the indole ring by alkyl, haloalkyl,  
6 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,  
7 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or  
8 2-(4-morpholinyl)ethyl whether or not further substituted  
9 in the indole ring to any extent, whether or not  
10 substituted in the naphthyl ring to any extent. Examples  
11 of this structural class include, but are not limited to,  
12 JWH-018, AM-2201, JWH-175, JWH-184, and JWH-185;

13 (43) Any compound structurally derived from  
14 3-(1-naphthoyl)pyrrole by substitution at the nitrogen  
15 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl,  
16 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
17 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
18 2-(4-morpholinyl)ethyl, whether or not further substituted  
19 in the pyrrole ring to any extent, whether or not  
20 substituted in the naphthyl ring to any extent. Examples  
21 of this structural class include, but are not limited to,  
22 JWH-030, JWH-145, JWH-146, JWH-307, and JWH-368;

23 (44) Any compound structurally derived from  
24 1-(1-naphthylmethyl)indene by substitution at the  
25 3-position of the indene ring by alkyl, haloalkyl,  
26 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,

1 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or  
2 2-(4-morpholinyl)ethyl whether or not further substituted  
3 in the indene ring to any extent, whether or not  
4 substituted in the naphthyl ring to any extent. Examples  
5 of this structural class include, but are not limited to,  
6 JWH-176;

7 (45) Any compound structurally derived from  
8 3-phenylacetylindole by substitution at the nitrogen atom  
9 of the indole ring with alkyl, haloalkyl, alkenyl,  
10 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
11 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
12 2-(4-morpholinyl)ethyl, whether or not further substituted  
13 in the indole ring to any extent, whether or not  
14 substituted in the phenyl ring to any extent. Examples of  
15 this structural class include, but are not limited to,  
16 JWH-167, JWH-250, JWH-251, and RCS-8;

17 (46) Any compound structurally derived from  
18 2-(3-hydroxycyclohexyl)phenol by substitution at the  
19 5-position of the phenolic ring by alkyl, haloalkyl,  
20 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,  
21 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or  
22 2-(4-morpholinyl)ethyl, whether or not substituted in the  
23 cyclohexyl ring to any extent. Examples of this structural  
24 class include, but are not limited to, CP 47, 497 and its  
25 C8 homologue (cannabicyclohexanol);

26 (46.1) Any compound structurally derived from

1 3-(benzoyl) indole with substitution at the nitrogen atom  
2 of the indole ring by an alkyl, haloalkyl, alkenyl,  
3 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
4 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
5 2-(4-morpholinyl)ethyl group whether or not further  
6 substituted in the indole ring to any extent and whether  
7 or not substituted in the phenyl ring to any extent.  
8 Examples of this structural class include, but are not  
9 limited to, AM-630, AM-2233, AM-694, Pravadoline (WIN  
10 48,098), and RCS-4;

11 (47) (Blank);

12 (48) (Blank);

13 (49) (Blank);

14 (50) (Blank);

15 (51) (Blank);

16 (52) (Blank);

17 (53) 2,5-Dimethoxy-4-(n)-propylthio-phenethylamine.

18 Some trade or other names: 2C-T-7;

19 (53.1) 4-ethyl-2,5-dimethoxyphenethylamine. Some  
20 trade or other names: 2C-E;

21 (53.2) 2,5-dimethoxy-4-methylphenethylamine. Some  
22 trade or other names: 2C-D;

23 (53.3) 4-chloro-2,5-dimethoxyphenethylamine. Some  
24 trade or other names: 2C-C;

25 (53.4) 4-iodo-2,5-dimethoxyphenethylamine. Some trade  
26 or other names: 2C-I;

1 (53.5) 4-ethylthio-2,5-dimethoxyphenethylamine. Some  
2 trade or other names: 2C-T-2;

3 (53.6) 2,5-dimethoxy-4-isopropylthio-phenethylamine.  
4 Some trade or other names: 2C-T-4;

5 (53.7) 2,5-dimethoxyphenethylamine. Some trade or  
6 other names: 2C-H;

7 (53.8) 2,5-dimethoxy-4-nitrophenethylamine. Some  
8 trade or other names: 2C-N;

9 (53.9) 2,5-dimethoxy-4-(n)-propylphenethylamine. Some  
10 trade or other names: 2C-P;

11 (53.10) 2,5-dimethoxy-3,4-dimethylphenethylamine.  
12 Some trade or other names: 2C-G;

13 (53.11) The N-(2-methoxybenzyl) derivative of any 2C  
14 phenethylamine referred to in subparagraphs (20.1), (53),  
15 (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7),  
16 (53.8), (53.9), and (53.10) including, but not limited to,  
17 25I-NBOMe and 25C-NBOMe;

18 (54) 5-Methoxy-N,N-diisopropyltryptamine;

19 (55) (Blank);

20 (56) (Blank);

21 (57) (Blank);

22 (58) (Blank);

23 (59) 3-cyclopropoylindole with substitution at the  
24 nitrogen atom of the indole ring by alkyl, haloalkyl,  
25 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,  
26 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or

1 2-(4-morpholinyl)ethyl, whether or not further substituted  
2 on the indole ring to any extent, whether or not  
3 substituted on the cyclopropyl ring to any extent:  
4 including, but not limited to, XLR11, UR144, FUB-144;

5 (60) 3-adamantoylindole with substitution at the  
6 nitrogen atom of the indole ring by alkyl, haloalkyl,  
7 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,  
8 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or  
9 2-(4-morpholinyl)ethyl, whether or not further substituted  
10 on the indole ring to any extent, whether or not  
11 substituted on the adamantyl ring to any extent:  
12 including, but not limited to, AB-001;

13 (61) N-(adamantyl)-indole-3-carboxamide with  
14 substitution at the nitrogen atom of the indole ring by  
15 alkyl, haloalkyl, alkenyl, cycloalkylmethyl,  
16 cycloalkylethyl, aryl halide, alkyl aryl halide,  
17 1-(N-methyl-2-piperidinyl)methyl, or  
18 2-(4-morpholinyl)ethyl, whether or not further substituted  
19 on the indole ring to any extent, whether or not  
20 substituted on the adamantyl ring to any extent:  
21 including, but not limited to, APICA/2NE-1, STS-135;

22 (62) N-(adamantyl)-indazole-3-carboxamide with  
23 substitution at a nitrogen atom of the indazole ring by  
24 alkyl, haloalkyl, alkenyl, cycloalkylmethyl,  
25 cycloalkylethyl, aryl halide, alkyl aryl halide,  
26 1-(N-methyl-2-piperidinyl)methyl, or

1 2-(4-morpholinyl)ethyl, whether or not further substituted  
2 on the indazole ring to any extent, whether or not  
3 substituted on the adamantyl ring to any extent:  
4 including, but not limited to, AKB48, 5F-AKB48;

5 (63) 1H-indole-3-carboxylic acid 8-quinolinyl ester  
6 with substitution at the nitrogen atom of the indole ring  
7 by alkyl, haloalkyl, alkenyl, cycloalkylmethyl,  
8 cycloalkylethyl, aryl halide, alkyl aryl halide,  
9 1-(N-methyl-2-piperidinyl)methyl, or  
10 2-(4-morpholinyl)ethyl, whether or not further substituted  
11 on the indole ring to any extent, whether or not  
12 substituted on the quinoline ring to any extent:  
13 including, but not limited to, PB22, 5F-PB22, FUB-PB-22;

14 (64) 3-(1-naphthoyl)indazole with substitution at the  
15 nitrogen atom of the indazole ring by alkyl, haloalkyl,  
16 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,  
17 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or  
18 2-(4-morpholinyl)ethyl, whether or not further substituted  
19 on the indazole ring to any extent, whether or not  
20 substituted on the naphthyl ring to any extent: including,  
21 but not limited to, THJ-018, THJ-2201;

22 (65) 2-(1-naphthoyl)benzimidazole with substitution  
23 at the nitrogen atom of the benzimidazole ring by alkyl,  
24 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,  
25 aryl halide, alkyl aryl halide,  
26 1-(N-methyl-2-piperidinyl)methyl, or



1 2-(4-morpholinyl)ethyl, whether or not further substituted  
2 on the benzimidazole ring to any extent, whether or not  
3 substituted on the naphthyl ring to any extent: including,  
4 but not limited to, FUBIMINA;

5 (66)

6 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indazole-  
7 3-carboxamide with substitution on the nitrogen atom of  
8 the indazole ring by alkyl, haloalkyl, alkenyl,  
9 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
10 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
11 2-(4-morpholinyl)ethyl, whether or not further substituted  
12 on the indazole ring to any extent: including, but not  
13 limited to, AB-PINACA, AB-FUBINACA, AB-CHMINACA;

14 (67) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-

15 indazole-3-carboxamide with substitution on the nitrogen  
16 atom of the indazole ring by alkyl, haloalkyl, alkenyl,  
17 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
18 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
19 2-(4-morpholinyl)ethyl, whether or not further substituted  
20 on the indazole ring to any extent: including, but not  
21 limited to, ADB-PINACA, ADB-FUBINACA;

22 (68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-

23 indole-3-carboxamide with substitution on the nitrogen  
24 atom of the indole ring by alkyl, haloalkyl, alkenyl,  
25 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
26 halide, 1-(N-methyl-2-piperidinyl)methyl, or

1 2-(4-morpholinyl)ethyl, whether or not further substituted  
2 on the indole ring to any extent: including, but not  
3 limited to, ADBICA, 5F-ADBICA;

4 (69) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indole-  
5 3-carboxamide with substitution on the nitrogen atom of  
6 the indole ring by alkyl, haloalkyl, alkenyl,  
7 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
8 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
9 2-(4-morpholinyl)ethyl, whether or not further substituted  
10 on the indole ring to any extent: including, but not  
11 limited to, ABICA, 5F-ABICA;

12 (70) Methyl 2-(1H-indazole-3-carboxamido)-3-  
13 methylbutanoate with substitution on the nitrogen atom of  
14 the indazole ring by alkyl, haloalkyl, alkenyl,  
15 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
16 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
17 2-(4-morpholinyl)ethyl, whether or not further substituted  
18 on the indazole ring to any extent: including, but not  
19 limited to, AMB, 5F-AMB;

20 (71) Methyl 2-(1H-indazole-3-carboxamido)-3,3-  
21 dimethylbutanoate with substitution on the nitrogen atom  
22 of the indazole ring by alkyl, haloalkyl, alkenyl,  
23 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
24 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
25 2-(4-morpholinyl)ethyl, whether or not further substituted  
26 on the indazole ring to any extent: including, but not

1 limited to, 5-fluoro-MDMB-PINACA, MDMB-FUBINACA;

2 (72) Methyl 2-(1H-indole-3-carboxamido)-3-  
3 methylbutanoate with substitution on the nitrogen atom of  
4 the indole ring by alkyl, haloalkyl, alkenyl,  
5 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
6 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
7 2-(4-morpholinyl)ethyl, whether or not further substituted  
8 on the indazole ring to any extent: including, but not  
9 limited to, MMB018, MMB2201, and AMB-CHMICA;

10 (73) Methyl 2-(1H-indole-3-carboxamido)-3,3-  
11 dimethylbutanoate with substitution on the nitrogen atom  
12 of the indole ring by alkyl, haloalkyl, alkenyl,  
13 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
14 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
15 2-(4-morpholinyl)ethyl, whether or not further substituted  
16 on the indazole ring to any extent: including, but not  
17 limited to, MDMB-CHMICA;

18 (74) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-  
19 indazole-3-carboxamide with substitution on the nitrogen  
20 atom of the indazole ring by alkyl, haloalkyl, alkenyl,  
21 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
22 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
23 2-(4-morpholinyl)ethyl, whether or not further substituted  
24 on the indazole ring to any extent: including, but not  
25 limited to, APP-CHMINACA, 5-fluoro-APP-PINACA;

26 (75) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-indole-

1 3-carboxamide with substitution on the nitrogen atom of  
2 the indole ring by alkyl, haloalkyl, alkenyl,  
3 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl  
4 halide, 1-(N-methyl-2-piperidinyl)methyl, or  
5 2-(4-morpholinyl)ethyl, whether or not further substituted  
6 on the indazole ring to any extent: including, but not  
7 limited to, APP-PICA and 5-fluoro-APP-PICA;

8 (76) 4-Acetoxy-N,N-dimethyltryptamine: trade name  
9 4-AcO-DMT;

10 (77) 5-Methoxy-N-methyl-N-isopropyltryptamine: trade  
11 name 5-MeO-MIPT;

12 (78) 4-hydroxy Diethyltryptamine (4-HO-DET);

13 (79) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);

14 (80) 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT);

15 (81) 4-hydroxy-N-methyl-N-isopropyltryptamine  
16 (4-HO-MiPT);

17 (82) Fluorophenylpiperazine;

18 (83) Methoxetamine;

19 (84) 1-(Ethylamino)-2-phenylpropan-2-one (iso-  
20 ethcathinone).

21 (e) Unless specifically excepted or unless listed in  
22 another schedule, any material, compound, mixture, or  
23 preparation which contains any quantity of the following  
24 substances having a depressant effect on the central nervous  
25 system, including its salts, isomers, and salts of isomers  
26 whenever the existence of such salts, isomers, and salts of

1 isomers is possible within the specific chemical designation:

2 (1) mecloqualone;

3 (2) methaqualone; and

4 (3) gamma hydroxybutyric acid.

5 (f) Unless specifically excepted or unless listed in  
6 another schedule, any material, compound, mixture, or  
7 preparation which contains any quantity of the following  
8 substances having a stimulant effect on the central nervous  
9 system, including its salts, isomers, and salts of isomers:

10 (1) Fenethylline;

11 (2) N-ethylamphetamine;

12 (3) Aminorex (some other names:

13 2-amino-5-phenyl-2-oxazoline; aminoxaphen;

14 4-5-dihydro-5-phenyl-2-oxazolamine) and its

15 salts, optical isomers, and salts of optical isomers;

16 (4) Methcathinone (some other names:

17 2-methylamino-1-phenylpropan-1-one;

18 Ephedrone; 2-(methylamino)-propiofenone;

19 alpha-(methylamino)propiofenone; N-methylcathinone;

20 methcathinone; Monomethylpropion; UR 1431) and its

21 salts, optical isomers, and salts of optical isomers;

22 (5) Cathinone (some trade or other names:

23 2-aminopropiofenone; alpha-aminopropiofenone;

24 2-amino-1-phenyl-propanone; norephedrone);

25 (6) N,N-dimethylamphetamine (also known as:

26 N,N-alpha-trimethyl-benzeneethanamine;

1 N,N-alpha-trimethylphenethylamine);

2 (7) (+ or -) cis-4-methylaminorex ((+ or -) cis-  
3 4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);

4 (8) 3,4-Methylenedioxypropylamphetamine (MDPV);

5 (9) Halogenated amphetamines and  
6 methamphetamines - any compound derived from either  
7 amphetamine or methamphetamine through the substitution  
8 of a halogen on the phenyl ring, including, but not  
9 limited to, 2-fluoroamphetamine, 3-  
10 fluoroamphetamine and 4-fluoroamphetamine;

11 (10) Aminopropylbenzofuran (APB):  
12 including 4-(2-Aminopropyl) benzofuran, 5-  
13 (2-Aminopropyl)benzofuran, 6-(2-Aminopropyl)  
14 benzofuran, and 7-(2-Aminopropyl) benzofuran;

15 (11) Aminopropyl-dihydrobenzofuran (APDB):  
16 including 4-(2-Aminopropyl)-2,3- dihydrobenzofuran,  
17 5-(2-Aminopropyl)-2, 3-dihydrobenzofuran,  
18 6-(2-Aminopropyl)-2,3-dihydrobenzofuran,  
19 and 7-(2-Aminopropyl)-2,3-dihydrobenzofuran;

20 (12) Methylaminopropylbenzofuran  
21 (MAPB): including 4-(2-methylaminopropyl)  
22 benzofuran, 5-(2-methylaminopropyl)benzofuran,  
23 6-(2-methylaminopropyl)benzofuran  
24 and 7-(2-methylaminopropyl)benzofuran.

25 (g) Temporary listing of substances subject to emergency  
26 scheduling. Any material, compound, mixture, or preparation

1 that contains any quantity of the following substances:

2 (1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide  
3 (benzylfentanyl), its optical isomers, isomers, salts, and  
4 salts of isomers;

5 (2) N-[1(2-thienyl) methyl-4-piperidyl]-N-  
6 phenylpropanamide (thenylfentanyl), its optical isomers,  
7 salts, and salts of isomers.

8 (h) Synthetic cathinones. Unless specifically excepted,  
9 any chemical compound which is not approved by the United  
10 States Food and Drug Administration or, if approved, is not  
11 dispensed or possessed in accordance with State or federal  
12 law, not including bupropion, structurally derived from  
13 2-aminopropan-1-one by substitution at the 1-position with  
14 either phenyl, naphthyl, or thiophene ring systems, whether or  
15 not the compound is further modified in one or more of the  
16 following ways:

17 (1) by substitution in the ring system to any extent  
18 with alkyl, alkylenedioxy, alkoxy, haloalkyl, hydroxyl, or  
19 halide substituents, whether or not further substituted in  
20 the ring system by one or more other univalent  
21 substituents. Examples of this class include, but are not  
22 limited to, 3,4-Methylenedioxycathinone (bk-MDA);

23 (2) by substitution at the 3-position with an acyclic  
24 alkyl substituent. Examples of this class include, but are  
25 not limited to, 2-methylamino-1-phenylbutan-1-one  
26 (buphedrone); or

1 (3) by substitution at the 2-amino nitrogen atom with  
2 alkyl, dialkyl, benzyl, or methoxybenzyl groups, or by  
3 inclusion of the 2-amino nitrogen atom in a cyclic  
4 structure. Examples of this class include, but are not  
5 limited to, Dimethylcathinone, Ethcathinone, and  
6 a-Pyrrolidinopropiophenone (a-PPP); or

7 Any other synthetic cathinone which is not approved by the  
8 United States Food and Drug Administration or, if approved, is  
9 not dispensed or possessed in accordance with State or federal  
10 law.

11 (i) Synthetic cannabinoids or piperazines. Any synthetic  
12 cannabinoid or piperazine which is not approved by the United  
13 States Food and Drug Administration or, if approved, which is  
14 not dispensed or possessed in accordance with State and  
15 federal law.

16 (j) Unless specifically excepted or listed in another  
17 schedule, any chemical compound which is not approved by the  
18 United States Food and Drug Administration or, if approved, is  
19 not dispensed or possessed in accordance with State or federal  
20 law, and is derived from the following structural classes and  
21 their salts:

22 (1) Benzodiazepine class: A fused 1,4-diazepine and  
23 benzene ring structure with a phenyl connected to the  
24 1,4-diazepine ring, with any substitution(s) or  
25 replacement(s) on the 1,4-diazepine or benzene ring, any  
26 substitution(s) on the phenyl ring, or any combination



1 thereof. Examples of this class include but are not  
2 limited to: Clonazolam, Flualprazolam; or

3 (2) Thienodiazepine class: A fused 1,4-diazepine and  
4 thiophene ring structure with a phenyl connected to the  
5 1,4-diazepine ring, with any substitution(s) or  
6 replacement(s) on the 1,4-diazepine or thiophene ring, any  
7 substitution(s) on the phenyl ring, or any combination  
8 thereof. Examples of this class include but are not  
9 limited to: Etizolam.

10 (Source: P.A. 99-371, eff. 1-1-16; 100-201, eff. 8-18-17;  
11 100-368, eff. 1-1-18; 100-789, eff. 1-1-19; 100-863, eff.  
12 8-14-18.)