

SB0702



100TH GENERAL ASSEMBLY

State of Illinois

2017 and 2018

SB0702

Introduced 1/30/2017, by Sen. Kyle McCarter

SYNOPSIS AS INTRODUCED:

720 ILCS 570/204

from Ch. 56 1/2, par. 1204

Amends the Illinois Controlled Substances Act. Adds 3,4-Dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (some trade or other name: U-47700) as a Schedule I controlled substance. Effective immediately.

LRB100 07354 RLC 17417 b

CORRECTIONAL
BUDGET AND
IMPACT NOTE ACT
MAY APPLY

A BILL FOR

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 their
11 isomers, esters, ethers, salts, and salts of isomers, esters,
12 and ethers, whenever the existence of such isomers, esters,
13 ethers and salts is possible within the specific chemical
14 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) Difenoxyin;
- 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 (c) Unless specifically excepted or unless listed in
18 another schedule, any of the following opium derivatives, its
19 salts, isomers and salts of isomers, whenever the existence of
20 such salts, isomers and salts of isomers is possible within the
21 specific chemical designation:

- 22 (1) Acetorphine;
- 23 (2) Acetyldihydrocodeine;
- 24 (3) Benzylmorphine;
- 25 (4) Codeine methylbromide;
- 26 (5) Codeine-N-Oxide;

- 1 (6) Cyprenorphine;
- 2 (7) Desomorphine;
- 3 (8) Diacetyldihydromorphine (Dihydroheroin);
- 4 (9) Dihydromorphine;
- 5 (10) Drotebanol;
- 6 (11) Etorphine (except hydrochloride salt);
- 7 (12) Heroin;
- 8 (13) Hydromorphanol;
- 9 (14) Methyldesorphine;
- 10 (15) Methyldihydromorphine;
- 11 (16) Morphine methylbromide;
- 12 (17) Morphine methylsulfonate;
- 13 (18) Morphine-N-Oxide;
- 14 (19) Myrophine;
- 15 (20) Nicocodeine;
- 16 (21) Nicomorphine;
- 17 (22) Normorphine;
- 18 (23) Pholcodine;
- 19 (24) Thebacon.

20 (d) Unless specifically excepted or unless listed in
21 another schedule, any material, compound, mixture, or
22 preparation which contains any quantity of the following
23 hallucinogenic substances, or which contains any of its salts,
24 isomers and salts of isomers, whenever the existence of such
25 salts, isomers, and salts of isomers is possible within the
26 specific chemical designation (for the purposes of this

1 paragraph only, the term "isomer" includes the optical,
2 position and geometric isomers):

3 (1) 3,4-methylenedioxyamphetamine

4 (alpha-methyl,3,4-methylenedioxyphenethylamine,
5 methylenedioxyamphetamine, MDA);

6 (1.1) Alpha-ethyltryptamine

7 (some trade or other names: etryptamine;
8 MONASE; alpha-ethyl-1H-indole-3-ethanamine;
9 3-(2-aminobutyl)indole; a-ET; and AET);

10 (2) 3,4-methylenedioxymethamphetamine (MDMA);

11 (2.1) 3,4-methylenedioxy-N-ethylamphetamine

12 (also known as: N-ethyl-alpha-methyl-
13 3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE,
14 and MDEA);

15 (2.2) N-Benzylpiperazine (BZP);

16 (2.2-1) Trifluoromethylphenylpiperazine (TFMPP);

17 (3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);

18 (4) 3,4,5-trimethoxyamphetamine (TMA);

19 (5) (Blank);

20 (6) Diethyltryptamine (DET);

21 (7) Dimethyltryptamine (DMT);

22 (7.1) 5-Methoxy-diallyltryptamine;

23 (8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);

24 (9) Ibogaine (some trade and other names:

25 7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-
26 6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b]

1 indole; Tabernanthe iboga);

2 (10) Lysergic acid diethylamide;

3 (10.1) Salvinorin A;

4 (10.5) Salvia divinorum (meaning all parts of the plant
5 presently classified botanically as Salvia divinorum,
6 whether growing or not, the seeds thereof, any extract from
7 any part of that plant, and every compound, manufacture,
8 salts, isomers, and salts of isomers whenever the existence
9 of such salts, isomers, and salts of isomers is possible
10 within the specific chemical designation, derivative,
11 mixture, or preparation of that plant, its seeds or
12 extracts);

13 (11) 3,4,5-trimethoxyphenethylamine (Mescaline);

14 (12) Peyote (meaning all parts of the plant presently
15 classified botanically as Lophophora williamsii Lemaire,
16 whether growing or not, the seeds thereof, any extract from
17 any part of that plant, and every compound, manufacture,
18 salts, derivative, mixture, or preparation of that plant,
19 its seeds or extracts);

20 (13) N-ethyl-3-piperidyl benzilate (JB 318);

21 (14) N-methyl-3-piperidyl benzilate;

22 (14.1) N-hydroxy-3,4-methylenedioxyamphetamine

23 (also known as N-hydroxy-alpha-methyl-

24 3,4(methylenedioxy)phenethylamine and N-hydroxy MDA);

25 (15) Parahexyl; some trade or other names:

26 3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-

1 dibenzo (b,d) pyran; Synhexyl;
2 (16) Psilocybin;
3 (17) Psilocyn;
4 (18) Alpha-methyltryptamine (AMT);
5 (19) 2,5-dimethoxyamphetamine
6 (2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA);
7 (20) 4-bromo-2,5-dimethoxyamphetamine
8 (4-bromo-2,5-dimethoxy-alpha-methylphenethylamine;
9 4-bromo-2,5-DMA);
10 (20.1) 4-Bromo-2,5 dimethoxyphenethylamine.
11 Some trade or other names: 2-(4-bromo-
12 2,5-dimethoxyphenyl)-1-aminoethane;
13 alpha-desmethyl DOB, 2CB, Nexus;
14 (21) 4-methoxyamphetamine
15 (4-methoxy-alpha-methylphenethylamine;
16 paramethoxyamphetamine; PMA);
17 (22) (Blank);
18 (23) Ethylamine analog of phencyclidine.
19 Some trade or other names:
20 N-ethyl-1-phenylcyclohexylamine,
21 (1-phenylcyclohexyl) ethylamine,
22 N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE;
23 (24) Pyrrolidine analog of phencyclidine. Some trade
24 or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy,
25 PHP;
26 (25) 5-methoxy-3,4-methylenedioxy-amphetamine;

- 1 (26) 2,5-dimethoxy-4-ethylamphetamine
2 (another name: DOET);
- 3 (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine
4 (another name: TCPy);
- 5 (28) (Blank);
- 6 (29) Thiophene analog of phencyclidine (some trade
7 or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine;
8 2-thienyl analog of phencyclidine; TPCP; TCP);
- 9 (30) Bufotenine (some trade or other names:
10 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole;
11 3-(2-dimethylaminoethyl)-5-indolol;
12 5-hydroxy-N,N-dimethyltryptamine;
13 N,N-dimethylserotonin; mappine);
- 14 (31) 1-Pentyl-3-(1-naphthoyl)indole
15 Some trade or other names: JWH-018;
- 16 (32) 1-Butyl-3-(1-naphthoyl)indole
17 Some trade or other names: JWH-073;
- 18 (33) 1-[(5-fluoropentyl)-1H-indol-3-yl]-
19 (2-iodophenyl)methanone
20 Some trade or other names: AM-694;
- 21 (34) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
22 (2-methyloctan-2-yl)phenol
23 Some trade or other names: CP 47,497
24 and its C6, C8 and C9 homologs;
- 25 (34.5) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
26 (2-methyloctan-2-yl)phenol), where side chain n=5;

1 and homologues where side chain n=4, 6, or 7; Some
2 trade or other names: CP 47,497;

3 (35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-
4 (2-methyloctan-2-yl)-6a,7,
5 10,10a-tetrahydrobenzo[c]chromen-1-ol

6 Some trade or other names: HU-210;

7 (35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-
8 dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
9 tetrahydrobenzo[c]chromen-1-ol, its isomers,
10 salts, and salts of isomers; Some trade or other
11 names: HU-210, Dexanabinol;

12 (36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-
13 6,6-dimethyl-3-(2-methyloctan-2-yl)-
14 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol
15 Some trade or other names: HU-211;

16 (37) (2-methyl-1-propyl-1H-indol-
17 3-yl)-1-naphthalenyl-methanone
18 Some trade or other names: JWH-015;

19 (38) 4-methoxynaphthalen-1-yl-
20 (1-pentylindol-3-yl)methanone
21 Some trade or other names: JWH-081;

22 (39) 1-Pentyl-3-(4-methyl-1-naphthoyl)indole
23 Some trade or other names: JWH-122;

24 (40) 2-(2-methylphenyl)-1-(1-pentyl-
25 1H-indol-3-yl)-ethanone
26 Some trade or other names: JWH-251;

1 (41) 1-(2-cyclohexylethyl)-3-

2 (2-methoxyphenylacetyl)indole

3 Some trade or other names: RCS-8, BTW-8 and SR-18;

4 (42) Any compound structurally derived from

5 3-(1-naphthoyl)indole or 1H-indol-3-yl-

6 (1-naphthyl)methane by substitution at the

7 nitrogen atom of the indole ring by alkyl, haloalkyl,

8 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,

9 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl,

10 or 2-(4-morpholinyl)ethyl whether or not further

11 substituted in the indole ring to any extent, whether

12 or not substituted in the naphthyl ring to any extent.

13 Examples of this structural class include, but are

14 not limited to, JWH-018, AM-2201, JWH-175, JWH-184,

15 and JWH-185;

16 (43) Any compound structurally derived from

17 3-(1-naphthoyl)pyrrole by substitution at the nitrogen

18 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl,

19 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl

20 aryl halide, 1-(N-methyl-2-piperidinyl)methyl,

21 or 2-(4-morpholinyl)ethyl, whether or not further

22 substituted in the pyrrole ring to any extent, whether

23 or not substituted in the naphthyl ring to any extent.

24 Examples of this structural class include, but are not

25 limited to, JWH-030, JWH-145, JWH-146, JWH-307, and

26 JWH-368;

1 (44) Any compound structurally derived from
2 1-(1-naphthylmethyl)indene by substitution
3 at the 3-position of the indene ring by alkyl, haloalkyl,
4 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl
5 halide, alkyl aryl halide, 1-(N-methyl-
6 2-piperidinyl)methyl, or 2-(4-
7 morpholinyl)ethyl whether or not further substituted in
8 the indene ring to any extent, whether or not substituted
9 in the naphthyl ring to any extent. Examples of
10 this structural class include, but are not
11 limited to, JWH-176;

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

23 (46) Any compound structurally derived from
24 2-(3-hydroxycyclohexyl)phenol by substitution
25 at the 5-position of the phenolic ring by alkyl,
26 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,

1 aryl halide, alkyl aryl halide, 1-(N-methyl-2-
2 piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
3 whether or not substituted in the cyclohexyl ring to any
4 extent. Examples of this structural class
5 include, but are not limited to, CP 47,
6 497 and its C8 homologue (cannabicyclohexanol);

7 (46.1) Benzoylindoles: Any compound
8 containing a 3-(benzoyl) indole structure with
9 substitution at the nitrogen atom of the
10 indole ring by an alkyl, haloalkyl, alkenyl,
11 cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2-piperidinyl)methyl,
13 or 2-(4-morpholinyl)ethyl group
14 whether or not further substituted
15 in the indole ring to any extent and
16 whether or not substituted in the phenyl ring
17 to any extent. Examples of this structural class
18 include, but are not limited, to, AM-630,
19 AM-2233, AM-694, Pravadoline (WIN 48,098), and RCS-4;

20 (47) 3,4-Methylenedioxymethcathinone
21 Some trade or other names: Methylone;

22 (48) 3,4-Methyenedioxypyrovalerone
23 Some trade or other names: MDPV;

24 (49) 4-Methylmethcathinone
25 Some trade or other names: Mephedrone;

26 (50) 4-methoxymethcathinone;

- 1 (51) 4-Fluoromethcathinone;
- 2 (52) 3-Fluoromethcathinone;
- 3 (53) 2,5-Dimethoxy-4-(n)-propylthio-
- 4 phenethylamine;
- 5 (54) 5-Methoxy-N,N-diisopropyltryptamine;
- 6 (55) Pentedrone;
- 7 (56) 4-iodo-2,5-dimethoxy-N-(2-methoxy
- 8 phenyl)methyl)-benzeneethanamine
- 9 (trade or other name: 25I-NBOMe);
- 10 (57) 4-chloro-2,5-dimethoxy-N-(2-methoxyphenyl)
- 11 methyl)-benzeneethanamine (trade or other name:
- 12 25C-NBOMe);
- 13 (58) 4-bromo-2,5-dimethoxy-N-(2-methoxyphenyl)
- 14 methyl)-benzeneethanamine (trade or other name:
- 15 25B-NBOMe);
- 16 (59) 3-cyclopropoylindole with
- 17 substitution at the nitrogen atom of the
- 18 indole ring by alkyl, haloalkyl, alkenyl,
- 19 cycloalkylmethyl, cycloalkylethyl, aryl
- 20 halide, alkyl aryl halide,
- 21 1-(N-methyl-2-piperidinyl)methyl, or
- 22 2-(4-morpholinyl)ethyl, whether or not
- 23 further substituted on the indole ring
- 24 to any extent, whether or not substituted
- 25 on the cyclopropyl ring to any extent:
- 26 including, but not limited to, XLR11,

1 UR144, FUB-144;

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

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

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

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

7 (63) 1H-indole-3-carboxylic acid 8-quinolinyl
8 ester with substitution at the nitrogen atom of the
9 indole ring by alkyl, haloalkyl, alkenyl,
10 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl
11 aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or
12 2-(4-morpholinyl)ethyl, whether or not further
13 substituted on the indole ring to any extent,
14 whether or not substituted on the quinoline ring
15 to any extent: including, but not limited to, PB22,
16 5F-PB22, FUB-PB-22;

17 (64) 3-(1-naphthoyl)indazole with
18 substitution at the nitrogen atom of the
19 indazole ring by alkyl, haloalkyl,
20 alkenyl, cycloalkylmethyl, cycloalkylethyl,
21 aryl halide, alkyl aryl halide,
22 1-(N-methyl-2-piperidinyl)methyl, or
23 2-(4-morpholinyl)ethyl, whether or not further
24 substituted on the indazole ring to any extent,
25 whether or not substituted on the naphthyl ring
26 to any extent: including, but not limited to,

1 THJ-018, THJ-2201;

2 (65) 2-(1-naphthoyl)benzimidazole with
3 substitution at the nitrogen atom of the benzimidazole
4 ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
5 cycloalkylethyl, aryl halide, alkyl aryl halide,
6 1-(N-methyl-2-piperidinyl)methyl, or
7 2-(4-morpholinyl)ethyl, whether or not further
8 substituted on the benzimidazole ring to any extent,
9 whether or not substituted on the naphthyl ring to
10 any extent: including, but not limited to, FUBIMINA;

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

20 (67) N-(1-amino-3,3-dimethyl-1-oxobutan-
21 2-yl)-1H-indazole-3-carboxamide with substitution
22 on the nitrogen atom of the indazole ring by alkyl,
23 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
24 aryl halide, alkyl aryl halide, 1-(N-methyl-2-
25 piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether
26 or not further substituted on the indazole ring to any

1 extent: including, but not limited to,

2 ADB-PINACA, ADB-FUBINACA;

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

10 including, but not limited to, ADBICA, 5F-ADBICA;

11 (69) N-(1-amino-3-methyl-1-oxobutan-2-yl)-
12 1H-indole-3-carboxamide with substitution on the
13 nitrogen atom of the indole ring by alkyl, haloalkyl,
14 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl
15 halide, alkyl aryl halide, 1-(N-methyl-2-
16 piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
17 whether or not further substituted on the indole
18 ring to any extent: including, but not limited
19 to, ABICA, 5F-ABICA;

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

1 ring to any extent: including, but not
2 limited to, AMB, 5F-AMB;
3 (71) 3,4-Dichloro-N-[2-(dimethylamino)
4 cyclohexyl]-N-methylbenzamide (Some trade or other
5 name: U-47700).

6 (e) Unless specifically excepted or unless listed in
7 another schedule, any material, compound, mixture, or
8 preparation which contains any quantity of the following
9 substances having a depressant effect on the central nervous
10 system, including its salts, isomers, and salts of isomers
11 whenever the existence of such salts, isomers, and salts of
12 isomers is possible within the specific chemical designation:

- 13 (1) mecloqualone;
14 (2) methaqualone; and
15 (3) gamma hydroxybutyric acid.

16 (f) Unless specifically excepted or unless listed in
17 another schedule, any material, compound, mixture, or
18 preparation which contains any quantity of the following
19 substances having a stimulant effect on the central nervous
20 system, including its salts, isomers, and salts of isomers:

- 21 (1) Fenethylline;
22 (2) N-ethylamphetamine;
23 (3) Aminorex (some other names:
24 2-amino-5-phenyl-2-oxazoline; aminoxaphen;
25 4-5-dihydro-5-phenyl-2-oxazolamine) and its
26 salts, optical isomers, and salts of optical isomers;

1 (4) Methcathinone (some other names:
2 2-methylamino-1-phenylpropan-1-one;
3 Ephedrone; 2-(methylamino)-propiofenone;
4 alpha-(methylamino)propiofenone; N-methylcathinone;
5 methycathinone; Monomethylpropion; UR 1431) and its
6 salts, optical isomers, and salts of optical isomers;

7 (5) Cathinone (some trade or other names:
8 2-aminopropiofenone; alpha-aminopropiofenone;
9 2-amino-1-phenyl-propanone; norephedrone);

10 (6) N,N-dimethylamphetamine (also known as:
11 N,N-alpha-trimethyl-benzeneethanamine;
12 N,N-alpha-trimethylphenethylamine);

13 (7) (+ or -) cis-4-methylaminorex ((+ or -) cis-
14 4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);

15 (8) 3,4-Methylenedioxypropylamphetamine (MDPV).

16 (g) Temporary listing of substances subject to emergency
17 scheduling. Any material, compound, mixture, or preparation
18 that contains any quantity of the following substances:

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

22 (2) N-[1(2-thienyl)
23 methyl-4-piperidyl]-N-phenylpropanamide (thienylfentanyl),
24 its optical isomers, salts, and salts of isomers.

25 (h) Synthetic cathinones. Unless specifically excepted,
26 any chemical compound not including bupropion, structurally

1 derived from 2-aminopropan-1-one by substitution at the
2 1-position with either phenyl, naphthyl, or thiophene ring
3 systems, whether or not the compound is further modified in one
4 or more of the following ways:

5 (1) by substitution in the ring system to
6 any extent with alkyl, alkylendioxy, alkoxy,
7 haloalkyl, hydroxyl, or halide substituents, whether
8 or not further substituted in the ring system
9 by one or more other univalent substituents.

10 Examples of this class include, but are not
11 limited to, 3,4-Methylenedioxycathinone
12 (bk-MDA);

13 (2) by substitution at the 3-position
14 with an acyclic alkyl substituent. Examples of
15 this class include, but are not limited to,
16 2-methylamino-1-phenylbutan-1-one
17 (buphedrone); or

18 (3) by substitution at the 2-amino nitrogen
19 atom with alkyl, dialkyl, benzyl, or methoxybenzyl
20 groups, or by inclusion of the 2-amino nitrogen atom
21 in a cyclic structure. Examples of this class include,
22 but are not limited to, Dimethylcathinone, Ethcathinone,
23 and α -Pyrrolidinopropiophenone (α -PPP).

24 (Source: P.A. 98-987, eff. 1-1-15; 99-371, eff. 1-1-16; revised
25 10-25-16.)

26 Section 99. Effective date. This Act takes effect upon

1 becoming law.