

100TH GENERAL ASSEMBLY

State of Illinois

2017 and 2018

HB3680

by Rep. Brandon W. Phelps

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: pink; U-47700) as a Schedule I controlled substance.

LRB100 10855 RLC 21089 b

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)

Sec. 204. (a) The controlled substances listed in this
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) Difenoxin;
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) Hydroxpethidine;
9	(28) Ketobemidone;
10	(29) Levomoramide;
11	(30) Levophenacylmorphan;
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	<pre>(3) Benzylmorphine;</pre>
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) Hydromorphinol;
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 list

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

- 6 - LRB100 10855 RLC 21089 b HB3680 paragraph only, the term "isomer" includes the optical, 1 2 position and geometric isomers): 3 (1) 3,4-methylenedioxyamphetamine 4 (alpha-methyl, 3, 4-methylenedioxyphenethylamine, 5 methylenedioxyamphetamine, MDA); (1.1) Alpha-ethyltryptamine 6 (some trade or other names: etryptamine; 7 MONASE; alpha-ethyl-1H-indole-3-ethanamine; 8 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, and MDEA); 14 15 (2.2) N-Benzylpiperazine (BZP); 16 (2.2-1) Trifluoromethylphenylpiperazine (TFMPP); 17 (3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA); (4) 3, 4, 5-trimethoxyamphetamine (TMA); 18 19 (5) (Blank); 20 (6) Diethyltryptamine (DET); (7) Dimethyltryptamine (DMT); 21 22 (7.1) 5-Methoxy-diallyltryptamine; 23 (8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP); (9) Ibogaine (some trade and other names: 24 25 7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b] 26

indole; Tabernanthe iboga);

2

(10) Lysergic acid diethylamide;

3

(10.1) Salvinorin A;

(10.5) Salvia divinorum (meaning all parts of the plant 4 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, salts, isomers, and salts of isomers whenever the existence 8 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);

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

20

21

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

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

(14.1) N-hydroxy-3,4-methylenedioxyamphetamine
(also known as N-hydroxy-alpha-methyl3,4 (methylenedioxy)phenethylamine and N-hydroxy MDA);

(15) Parahexyl; some trade or other names:
3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-

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HB3680
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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	<pre>paramethoxyamphetamine; PMA);</pre>
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;

- 9 -LRB100 10855 RLC 21089 b HB3680 (26) 2,5-dimethoxy-4-ethylamphetamine 1 2 (another name: DOET); (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine 3 (another name: TCPy); 4 5 (28) (Blank); (29) Thiophene analog of phencyclidine (some trade 6 7 or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 2-thienyl analog of phencyclidine; TPCP; TCP); 8 (30) Bufotenine (some trade or other names: 9 10 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole; 11 3-(2-dimethylaminoethyl)-5-indolol; 12 5-hydroxy-N,N-dimethyltryptamine; 13 N,N-dimethylserotonin; mappine); (31) 1-Pentyl-3-(1-naphthoyl)indole 14 15 Some trade or other names: JWH-018; 16 (32) 1-Butyl-3-(1-naphthoyl) indole 17 Some trade or other names: JWH-073; (33) 1-[(5-fluoropentyl)-1H-indol-3-yl]-18 (2-iodophenyl)methanone 19 Some trade or other names: AM-694; 20 21 (34) 2-[(1R, 3S)-3-hydroxycyclohexyl]-5-22 (2-methyloctan-2-yl)phenol 23 Some trade or other names: CP 47,497 and its C6, C8 and C9 homologs; 24 25 (34.5) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-26 (2-methyloctan-2-yl) phenol), where side chain n=5;

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and homologues where side chain n=4, 6, or 7; Some
1
 2
          trade or other names: CP 47,497;
               (35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-
 3
          (2-methyloctan-2-yl)-6a,7,
 4
 5
          10,10a-tetrahydrobenzo[c]chromen-1-ol
          Some trade or other names: HU-210;
 6
7
               (35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-
          dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
 8
 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)-
          6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol
14
15
          Some trade or other names: HU-211;
               (37) (2-methyl-1-propyl-1H-indol-
16
17
          3-yl)-1-naphthalenyl-methanone
          Some trade or other names: JWH-015;
18
               (38) 4-methoxynaphthalen-1-yl-
19
20
          (1-pentylindol-3-yl)methanone
          Some trade or other names: JWH-081;
21
22
               (39) 1-Pentyl-3-(4-methyl-1-naphthoyl) indole
23
          Some trade or other names: JWH-122;
               (40) 2-(2-methylphenyl)-1-(1-pentyl-
24
25
          1H-indol-3-yl)-ethanone
26
          Some trade or other names: JWH-251;
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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, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl 19 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 limited to, JWH-030, JWH-145, JWH-146, JWH-307, and 25 26 JWH-368;

- 12 - LRB100 10855 RLC 21089 b

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,

- 13 - LRB100 10855 RLC 21089 b

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 $_{\boldsymbol{L}}$ 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;

	HB3680	- 14	-	LRB100 1(0855 1	RLC 21089 b
1	(51)	4-Fluoromethcathin	one;			
2	(52)	3-Fluoromethcathin	one;			
3	(53)	2,5-Dimethoxy-4-(n)-propy	ylthio-		
4	phenethyl	amine;				
5	(54)	5-Methoxy-N,N-diis	opropyl	ltryptami	.ne;	
6	(55)	Pentedrone;				
7	(56)	4-iodo-2,5-dimetho	xy-N- ((2-methox	хy	
8	phenyl)me	thyl)-benzeneethana	mine			
9	(trade or	other name: 25I-NB	OMe);			
10	(57)	4-chloro-2,5-dimet	hoxy-N-	-[(2-meth	loxyp	henyl)
11	methyl]-b	enzeneethanamine (t	rade oi	r other n	ame:	
12	25C-NBOMe);				
13	(58)	4-bromo-2,5-dimeth	oxy-N-	[(2-methc	xyph	enyl)
14	methyl]-b	enzeneethanamine (t	rade oi	r other n	ame:	
15	25B-NBOMe);				
16	(59)	3-cyclopropoylindo	le with	ר		
17	substitut	ion at the nitrogen	atom o	of the		
18	indole ri	ng by alkyl, haloal	kyl, al	lkenyl,		
19	cycloalky	lmethyl, cycloalkyl	ethyl,	aryl		
20	halide, a	lkyl aryl halide,				
21	1-(N-meth	yl-2-piperidinyl)me	thyl, d	or		
22	2-(4-morp	nolinyl)ethyl, whet	her or	not		
23	further s	ubstituted on the is	ndole 1	ring		
24	to any ex	cent, whether or no	t subst	tituted		
25	on the cy	clopropyl ring to a	ny exte	ent:		
26	including	, but not limited t	o <u>,</u> XLR1	11,		

1 UR144, FUB-144;

26

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 <u>,</u>
12	but not limited to <u>,</u> 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 <u>,</u> but not limited to <u>,</u>
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,

cycloalkylethyl, aryl halide, alkyl aryl halide,

- 16 - LRB100 10855 RLC 21089 b

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, whether or not substituted on the adamantyl 4 5 ring to any extent: including, but not limited to, AKB48, 5F-AKB48; 6 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 substituted on the indole ring to any extent, 13 14 whether or not substituted on the guinoline ring 15 to any extent: including, but not limited to, PB22, 16 5F-PB22, FUB-PB-22; (64) 3-(1-naphthoyl) indazole with 17 18 substitution at the nitrogen atom of the 19 indazole ring by alkyl, haloalkyl, 20 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 21 22 1-(N-methyl-2-piperidinyl)methyl, or 23 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent, 24 25 whether or not substituted on the naphthyl ring 26 to any extent: including, but not limited to,

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1 THJ-018, THJ-2201;
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(65) 2-(1-naphthoyl)benzimidazole with
substitution at the nitrogen atom of the benzimidazole
ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
cycloalkylethyl, aryl halide, alkyl aryl halide,
1-(N-methyl-2-piperidinyl)methyl, or
2-(4-morpholinyl)ethyl, whether or not further
substituted on the benzimidazole ring to any extent,
whether or not substituted on the naphthyl ring to
any extent: including, but not limited to <u>,</u> FUBIMINA;
(66) N-(1-amino-3-methyl-1-oxobutan-2-yl)
-1H-indazole-3-carboxamide with substitution on the
nitrogen atom of the indazole ring by alkyl,
haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
aryl halide, alkyl aryl halide, 1-(N-methyl-2-
piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
whether or not further substituted on the indazole
ring to any extent: including $_{\it L}$ but not limited to $_{\it L}$
AB-PINACA, AB-FUBINACA, AB-CHMINACA;
(67) N-(1-amino-3,3-dimethyl-1-oxobutan-
2-yl)-1H-indazole-3-carboxamide with substitution
on the nitrogen atom of the indazole ring by alkyl,
haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
aryl halide, alkyl aryl halide, 1-(N-methyl-2-
piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether
or not further substituted on the indazole ring to any

- 18 - LRB100 10855 RLC 21089 b

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extent: including<u>,</u> but not limited to<u>,</u> ADB-PINACA, ADB-FUBINACA;

(68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-3 1H-indole-3-carboxamide with substitution on the nitrogen 4 5 atom of the indole ring by alkyl, haloalkyl, alkenyl, 6 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 7 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 nitrogen atom of the indole ring by alkyl, haloalkyl, 13 14 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl 15 halide, alkyl aryl halide, 1-(N-methyl-2piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, 16 17 whether or not further substituted on the indole ring to any extent: including, but not limited 18 19 to, ABICA, 5F-ABICA;

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

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1 ring to any extent: including, but not
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2 limited to, AMB, 5F-AMB;-

(71) 3,4-Dichloro-N-[2-(dimethylamino)

4 cyclohexyl]-N-methylbenzamide (Some trade or other

5 name: pink; U-47700).

HB3680

Unless specifically excepted or unless listed in 6 (e) another schedule, any material, compound, mixture, 7 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

3

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)-propiophenone;
4	alpha-(methylamino)propiophenone; 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-aminopropiophenone; alpha-aminopropiophenone;
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-Methylenedioxypyrovalerone (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	<pre>methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl),</pre>
24	its optical isomers, salts, and salts of isomers.
25	(h) Synthetic cathinones. Unless specifically excepted,
26	any chemical compound not including bupropion, structurally

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

5 (1) by substitution in the ring system to any extent with alkyl, alkylenedioxy, alkoxy, 6 7 haloalkyl, hydroxyl, or halide substituents, whether or not further substituted in the ring system 8 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);

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

17 (buphedrone); or

(3) by substitution at the 2-amino nitrogen 18 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 a-Pyrrolidinopropiophenone (a-PPP). (Source: P.A. 98-987, eff. 1-1-15; 99-371, eff. 1-1-16; revised 24 25 10 - 25 - 16.