

## 103RD GENERAL ASSEMBLY

# State of Illinois

# 2023 and 2024

### HB4352

Introduced 1/16/2024, by Rep. Tom Weber

## SYNOPSIS AS INTRODUCED:

720	ILCS	570/204	from (	Ch.	56	1/2,	par.	1204
720	ILCS	570/206	from (	Ch.	56	1/2,	par.	1206

Amends the Illinois Controlled Substances Act. Provides that Xylazine and Clonazolam are to be regulated under the Act as Schedule II controlled substances.

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AN ACT concerning criminal law.

# Be it enacted by the People of the State of Illinois, represented in the General Assembly:

4 Section 5. The Illinois Controlled Substances Act is 5 amended by changing Sections 204 and 206 as follows:

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

(Text of Section before amendment by P.A. 103-245)

8 Sec. 204. (a) The controlled substances listed in this 9 Section are included in Schedule I.

10 (b) Unless specifically excepted or unless listed in 11 another schedule, any of the following opiates, including 12 their isomers, esters, ethers, salts, and salts of isomers, 13 esters, and ethers, whenever the existence of such isomers, 14 esters, ethers and salts is possible within the specific 15 chemical designation:

16

20

(1) Acetylmethadol;

17 (1.1) Acetyl-alpha-methylfentanyl

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

19 4-piperidinyl]-N-phenylacetamide);

(2) Allylprodine;

21 (3) Alphacetylmethadol, except

22 levo-alphacetylmethadol (also known as levo-alpha-23 acetylmethadol, levomethadyl acetate, or LAAM);

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1	(4) Alphameprodine;						
2	(5) Alphamethadol;						
3	(6) Alpha-methylfenta	nyl					
4	(N-(1-alpha-methyl-beta-p	henyl)	) etl	nyl-4-j	piperidy	/l)	
5	propionanilide; 1-(1-met	hyl-2-	-phei	nyleth	yl)-4-(1	1—	
6	propanilido) piperidine;						
7	(6.1) Alpha-methylthi	ofenta	anyl				
8	(N-[1-methyl-2-(2-thienyl	)ethy]	1-				
9	4-piperidinyl]-N-phenylpr	opanan	mide	);			
10	(7) 1-methyl-4-phenyl-	-4-prc	opior	noxypip	peridine	) (I	MPPP);
11	(7.1) PEPAP						
12	(1-(2-phenethyl)-4-phenyl	-4-ace	etoxy	ypiper	idine);		
13	(8) Benzethidine;						
14	(9) Betacetylmethadol	÷					
15	(9.1) Beta-hydroxyfen	tanyl					
16	(N-[1-(2-hydroxy-2-phenet	hyl)-					
17	4-piperidinyl]-N-phenylpr	opanan	mide	;			
18	(10) Betameprodine;						
19	(11) Betamethadol;						
20	(12) Betaprodine;						
21	(13) Clonitazene;						
22	(14) Dextromoramide;						
23	(15) Diampromide;						
24	(16) Diethylthiambute	ne;					
25	(17) Difenoxin;						
26	(18) Dimenoxadol;						

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

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

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1	(4-nitrophenyl)ethyl]-2-piperidinylidene]-
2	benzenesulfonamide (W-18);
3	(56) 4-chloro-N-[1-(2-phenylethyl)
4	-2-piperidinylidene]-benzenesulfonamide (W-15);
5	(57) acrylfentanyl (acryloylfentanyl).
6	(c) Unless specifically excepted or unless listed in
7	another schedule, any of the following opium derivatives, its
8	salts, isomers and salts of isomers, whenever the existence of
9	such salts, isomers and salts of isomers is possible within
10	the specific chemical designation:
11	(1) Acetorphine;
12	(2) Acetyldihydrocodeine;
13	<pre>(3) Benzylmorphine;</pre>
14	(4) Codeine methylbromide;
15	(5) Codeine-N-Oxide;
16	(6) Cyprenorphine;
17	(7) Desomorphine;
18	(8) Diacetyldihydromorphine (Dihydroheroin);
19	(9) Dihydromorphine;
20	(10) Drotebanol;
21	(11) Etorphine (except hydrochloride salt);
22	(12) Heroin;
23	(13) Hydromorphinol;
24	(14) Methyldesorphine;
25	(15) Methyldihydromorphine;
26	(16) Morphine methylbromide;

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1	(17)	Morphine	<pre>methylsulfonate;</pre>

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

9 Unless specifically excepted or unless listed in (d) 10 another schedule, any material, compound, mixture, or 11 preparation which contains any quantity of the following 12 hallucinogenic substances, or which contains any of its salts, 13 isomers and salts of isomers, whenever the existence of such 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 17 position and geometric isomers):

- 18
- (1) 3,4-methylenedioxyamphetamine

19 (alpha-methyl, 3, 4-methylenedioxyphenethylamine,

- 20 methylenedioxyamphetamine, MDA);
- 21

(1.1) Alpha-ethyltryptamine

22 (some trade or other names: etryptamine;

- 23 MONASE; alpha-ethyl-1H-indole-3-ethanamine;
- 24 3-(2-aminobutyl)indole; a-ET; and AET);

25 (2) 3,4-methylenedioxymethamphetamine (MDMA);

26 (2.1) 3,4-methylenedioxy-N-ethylamphetamine

1	(also known as: N-ethyl-alpha-methyl-
2	3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE,
3	and MDEA);
4	(2.2) N-Benzylpiperazine (BZP);
5	(2.2-1) Trifluoromethylphenylpiperazine (TFMPP);
6	(3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);
7	(4) 3,4,5-trimethoxyamphetamine (TMA);
8	(5) (Blank);
9	(6) Diethyltryptamine (DET);
10	(7) Dimethyltryptamine (DMT);
11	(7.1) 5-Methoxy-diallyltryptamine;
12	<pre>(8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);</pre>
13	(9) Ibogaine (some trade and other names:
14	7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-
15	6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b]
16	<pre>indole; Tabernanthe iboga);</pre>
17	(10) Lysergic acid diethylamide;
18	(10.1) Salvinorin A;
19	(10.5) Salvia divinorum (meaning all parts of the
20	plant presently classified botanically as Salvia
21	divinorum, whether growing or not, the seeds thereof, any
22	extract from any part of that plant, and every compound,
23	manufacture, salts, isomers, and salts of isomers whenever
24	the existence of such salts, isomers, and salts of isomers
25	is possible within the specific chemical designation,
26	derivative, mixture, or preparation of that plant, its

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

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1	2,5-dimethoxyphenyl)-1-aminoethane;
2	alpha-desmethyl DOB, 2CB, Nexus;
3	(21) 4-methoxyamphetamine
4	(4-methoxy-alpha-methylphenethylamine;
5	<pre>paramethoxyamphetamine; PMA);</pre>
6	(22) (Blank);
7	(23) Ethylamine analog of phencyclidine.
8	Some trade or other names:
9	N-ethyl-1-phenylcyclohexylamine,
10	(1-phenylcyclohexyl) ethylamine,
11	N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE;
12	(24) Pyrrolidine analog of phencyclidine. Some trade
13	or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy,
14	PHP;
15	(25) 5-methoxy-3,4-methylenedioxy-amphetamine;
16	(26) 2,5-dimethoxy-4-ethylamphetamine
17	(another name: DOET);
18	(27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine
19	(another name: TCPy);
20	(28) (Blank);
21	(29) Thiophene analog of phencyclidine (some trade
22	or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine;
23	2-thienyl analog of phencyclidine; TPCP; TCP);
24	(29.1) Benzothiophene analog of phencyclidine. Some
25	trade or other names: BTCP or benocyclidine;
26	(29.2) 3-Methoxyphencyclidine (3-MeO-PCP);

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

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1 (40) (Blank);

(41) (Blank);

3 (42) Any compound structurally derived from 3-(1-naphthoyl) indole 4 or 5 1H-indol-3-yl-(1-naphthyl)methane by substitution at the 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 9 2-(4-morpholinyl)ethyl whether or not further substituted 10 in the indole ring to any extent, whether or not 11 substituted in the naphthyl ring to any extent. Examples 12 of this structural class include, but are not limited to, JWH-018, AM-2201, JWH-175, JWH-184, and JWH-185; 13

14 (43) Any compound structurally derived from 15 3-(1-naphthoyl)pyrrole by substitution at the nitrogen 16 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl, 17 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, 18 or 2-(4-morpholinyl)ethyl, whether or not further substituted 19 20 in the pyrrole ring to any extent, whether or not 21 substituted in the naphthyl ring to any extent. Examples 22 of this structural class include, but are not limited to, 23 JWH-030, JWH-145, JWH-146, JWH-307, and JWH-368;

(44) Any compound structurally derived from
1-(1-naphthylmethyl)indene by substitution at the
3-position of the indene 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 in the indene ring to any extent, whether or not substituted in the naphthyl ring to any extent. Examples of this structural class include, but are not limited to, JWH-176;

8 (45) compound structurally derived Any from 9 3-phenylacetylindole by substitution at the nitrogen atom 10 of the indole ring with alkyl, haloalkyl, alkenyl, 11 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 12 halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 13 14 in the indole ring to any extent, whether or not 15 substituted in the phenyl ring to any extent. Examples of 16 this structural class include, but are not limited to, 17 JWH-167, JWH-250, JWH-251, and RCS-8;

Any compound structurally derived 18 (46) from 19 2-(3-hydroxycyclohexyl)phenol by substitution at the 20 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, 21 22 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 23 2-(4-morpholinyl)ethyl, whether or not substituted in the cyclohexyl ring to any extent. Examples of this structural 24 25 class include, but are not limited to, CP 47, 497 and its 26 C8 homologue (cannabicyclohexanol);

1	(46.1) Any compound structurally derived from
2	3-(benzoyl) indole with substitution at the nitrogen atom
3	of the indole ring by an alkyl, haloalkyl, alkenyl,
4	cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl
5	halide, 1-(N-methyl-2-piperidinyl)methyl, or
6	2-(4-morpholinyl)ethyl group whether or not further
7	substituted in the indole ring to any extent and whether
8	or not substituted in the phenyl ring to any extent.
9	Examples of this structural class include, but are not
10	limited to, AM-630, AM-2233, AM-694, Pravadoline (WIN
11	48,098), and RCS-4;
12	(47) (Blank);
13	(48) (Blank);
14	(49) (Blank);
15	(50) (Blank);
16	(51) (Blank);
17	(52) (Blank);
18	(53) 2,5-Dimethoxy-4-(n)-propylthio-phenethylamine.
19	Some trade or other names: 2C-T-7;
20	(53.1) 4-ethyl-2,5-dimethoxyphenethylamine. Some
21	<pre>trade or other names: 2C-E;</pre>
22	(53.2) 2,5-dimethoxy-4-methylphenethylamine. Some
23	trade or other names: 2C-D;
24	(53.3) 4-chloro-2,5-dimethoxyphenethylamine. Some
25	trade or other names: 2C-C;
26	(53.4) 4-iodo-2,5-dimethoxyphenethylamine. Some trade

1	or other names: 2C-I;
2	(53.5) 4-ethylthio-2,5-dimethoxyphenethylamine. Some
3	trade or other names: 2C-T-2;
4	(53.6) 2,5-dimethoxy-4-isopropylthio-phenethylamine.
5	Some trade or other names: 2C-T-4;
6	(53.7) 2,5-dimethoxyphenethylamine. Some trade or
7	other names: 2C-H;
8	(53.8) 2,5-dimethoxy-4-nitrophenethylamine. Some
9	trade or other names: 2C-N;
10	(53.9) 2,5-dimethoxy-4-(n)-propylphenethylamine. Some
11	trade or other names: 2C-P;
12	(53.10) 2,5-dimethoxy-3,4-dimethylphenethylamine.
13	Some trade or other names: 2C-G;
1 /	(52.11) The N (2 methemoly device from 20
14	(53.11) The $N-(2-methoxypenzy1)$ derivative of any 20
15	phenethylamine referred to in subparagraphs (20.1), (53),
14 15 16	(53.11) The N-(2-methoxybenzyl) derivative of any 20 phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7),
14 15 16 17	(53.11) The N-(2-methoxybenzyl) derivative of any 20 phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to,
14 15 16 17 18	(53.11) The N-(2-Methoxybenzyl) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe;
14 15 16 17 18 19	<pre>(53.11) The N-(2-Methoxybenzy1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine;</pre>
14 15 16 17 18 19 20	<pre>(53.11) The N-(2-Methoxybenzy1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank);</pre>
14 15 16 17 18 19 20 21	<pre>(53.11) The N-(2-Methoxybenzy1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank); (56) (Blank);</pre>
14 15 16 17 18 19 20 21 22	<pre>(53.11) The N-(2-Methoxyben2y1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank); (56) (Blank); (57) (Blank);</pre>
14 15 16 17 18 19 20 21 22 23	<pre>(53.11) The N-(2-Methoxybenzy1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank); (56) (Blank); (57) (Blank); (58) (Blank);</pre>
14 15 16 17 18 19 20 21 22 23 24	<pre>(53.11) The N-(2-Methoxybenzy1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank); (56) (Blank); (57) (Blank); (58) (Blank); (59) 3-cyclopropoylindole with substitution at the</pre>
14 15 16 17 18 19 20 21 22 23 24 25	<pre>(53.11) The N-(2-Methoxyben2y1) derivative of any 2c phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe; (54) 5-Methoxy-N,N-diisopropyltryptamine; (55) (Blank); (56) (Blank); (57) (Blank); (58) (Blank); (59) 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl,</pre>

1 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2 2-(4-morpholinyl)ethyl, whether or not further substituted 3 on the indole ring to any extent, whether or not 4 substituted on the cyclopropyl ring to any extent: 5 including, but not limited to, XLR11, UR144, FUB-144;

(60) 3-adamantoylindole with substitution at 6 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, 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 adamantyl ring to any extent: including, but not limited to, AB-001; 13

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

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

or

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1-(N-methyl-2-piperidinyl)methyl,

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

(63) 1H-indole-3-carboxylic acid 8-quinolinyl ester 6 7 with substitution at the nitrogen atom of the indole ring 8 alkyl, haloalkyl, alkenyl, cycloalkylmethyl, by 9 cycloalkylethyl, aryl halide, alkyl aryl halide, 10 1-(N-methyl-2-piperidinyl)methyl, or 11 2-(4-morpholinyl)ethyl, whether or not further substituted 12 on the indole ring to any extent, whether or not the quinoline ring to 13 substituted on any extent: 14 including, but not limited to, PB22, 5F-PB22, FUB-PB-22;

15 (64) 3-(1-naphthoyl)indazole with substitution at the 16 nitrogen atom of the indazole ring by alkyl, haloalkyl, 17 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 18 2-(4-morpholinyl)ethyl, whether or not further substituted 19 20 on the indazole ring to any extent, whether or not substituted on the naphthyl ring to any extent: including, 21 22 but not limited to, THJ-018, THJ-2201;

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

or

1

1-(N-methyl-2-piperidinyl)methyl,

2 2-(4-morpholinyl)ethyl, whether or not further substituted 3 on the benzimidazole ring to any extent, whether or not 4 substituted on the naphthyl ring to any extent: including, 5 but not limited to, FUBIMINA;

6

(66)

7 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indazole-

3-carboxamide with substitution on the nitrogen atom of 8 9 indazole ring by alkyl, haloalkyl, the alkenvl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 10 11 halide, 1-(N-methyl-2-piperidinyl)methyl, or 12 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not 13 14 limited to, AB-PINACA, AB-FUBINACA, AB-CHMINACA;

15 (67)N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-16 indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, 17 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 18 1-(N-methyl-2-piperidinyl)methyl, 19 halide, or 20 2-(4-morpholinyl)ethyl, whether or not further substituted 21 on the indazole ring to any extent: including, but not 22 limited to, ADB-PINACA, ADB-FUBINACA;

(68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H indole-3-carboxamide with substitution on the nitrogen
 atom of the indole ring by alkyl, haloalkyl, alkenyl,
 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl

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

5 (69) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of 6 7 indole ring by alkyl, haloalkyl, alkenyl, the 8 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 9 halide, 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: including, but not 12 limited to, ABICA, 5F-ABICA;

2-(1H-indazole-3-carboxamido)-3-13 (70)Methyl 14 methylbutanoate with substitution on the nitrogen atom of indazole 15 the ring by alkyl, haloalkyl, alkenyl, 16 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 1-(N-methyl-2-piperidinyl)methyl, 17 halide, or 2-(4-morpholinyl)ethyl, whether or not further substituted 18 19 on the indazole ring to any extent: including, but not limited to, AMB, 5F-AMB; 20

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

1 2 on the indazole ring to any extent: including, but not limited to, 5-fluoro-MDMB-PINACA, MDMB-FUBINACA;

2-(1H-indole-3-carboxamido)-3-3 (72)Methyl methylbutanoate with substitution on the nitrogen atom of 4 5 the indole ring by alkyl, haloalkyl, alkenvl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 6 7 halide, 1-(N-methyl-2-piperidinyl)methyl, or 8 2-(4-morpholinyl)ethyl, whether or not further substituted 9 on the indazole ring to any extent: including, but not 10 limited to, MMB018, MMB2201, and AMB-CHMICA;

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

19 (74)N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1Hindazole-3-carboxamide with substitution on the nitrogen 20 21 atom of the indazole ring by alkyl, haloalkyl, alkenyl, 22 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 23 1-(N-methyl-2-piperidinyl)methyl, halide, or 24 2-(4-morpholinyl)ethyl, whether or not further substituted 25 on the indazole ring to any extent: including, but not limited to, APP-CHMINACA, 5-fluoro-APP-PINACA; 26

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(75) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-indole-1 2 3-carboxamide with substitution on the nitrogen atom of 3 the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 4 5 halide. 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 6 7 on the indazole ring to any extent: including, but not limited to, APP-PICA and 5-fluoro-APP-PICA; 8 9 (76) 4-Acetoxy-N, N-dimethyltryptamine: trade name 10 4 - ACO - DMT; 11 (77) 5-Methoxy-N-methyl-N-isopropyltryptamine: trade 12 name 5-MeO-MIPT; 13 (78) 4-hydroxy Diethyltryptamine (4-HO-DET); (79) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET); 14 15 (80) 4-hydroxy-N, N-diisopropyltryptamine (4-HO-DiPT); 16 (81) 4-hydroxy-N-methyl-N-isopropyltryptamine 17 (4-HO-MiPT); (82) Fluorophenylpiperazine; 18 (83) Methoxetamine; 19 20 (84)1-(Ethylamino)-2-phenylpropan-2-one (isoethcathinone). 21 22 (e) Unless specifically excepted or unless listed in 23 schedule, any material, compound, mixture, another or preparation which contains any quantity of the following 24 25 substances having a depressant effect on the central nervous 26 system, including its salts, isomers, and salts of isomers

whenever the existence of such salts, isomers, and salts of 1 2 isomers is possible within the specific chemical designation: 3 (1) mecloqualone; (2) methaqualone; and 4 5 (3) gamma hydroxybutyric acid. 6 (f) 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 stimulant effect on the central nervous 10 system, including its salts, isomers, and salts of isomers: 11 (1) Fenethylline; 12 (2) N-ethylamphetamine; 13 (3) Aminorex (some other names: 2-amino-5-phenyl-2-oxazoline; aminoxaphen; 14 15 4-5-dihydro-5-phenyl-2-oxazolamine) and its 16 salts, optical isomers, and salts of optical isomers; 17 (4) Methcathinone (some other names: 2-methylamino-1-phenylpropan-1-one; 18 19 Ephedrone; 2-(methylamino)-propiophenone; 20 alpha-(methylamino)propiophenone; N-methylcathinone; 21 methycathinone; Monomethylpropion; UR 1431) and its 22 salts, optical isomers, and salts of optical isomers; 23 (5) Cathinone (some trade or other names: 24 2-aminopropiophenone; alpha-aminopropiophenone; 25 2-amino-1-phenyl-propanone; norephedrone); 26 (6) N, N-dimethylamphetamine (also known as:

1	N,N-alpha-trimethyl-benzeneethanamine;
2	N,N-alpha-trimethylphenethylamine);
3	(7) (+ or -) cis-4-methylaminorex ((+ or -) cis-
4	4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);
5	<pre>(8) 3,4-Methylenedioxypyrovalerone (MDPV);</pre>
6	(9) Halogenated amphetamines and
7	methamphetamines - any compound derived from either
8	amphetamine or methamphetamine through the substitution
9	of a halogen on the phenyl ring, including, but not
10	limited to, 2-fluoroamphetamine, 3-
11	fluoroamphetamine and 4-fluoroamphetamine;
12	(10) Aminopropylbenzofuran (APB):
13	including 4-(2-Aminopropyl) benzofuran, 5-
14	(2-Aminopropyl)benzofuran, 6-(2-Aminopropyl)
15	benzofuran, and 7-(2-Aminopropyl) benzofuran;
16	(11) Aminopropyldihydrobenzofuran (APDB):
17	including 4-(2-Aminopropyl)-2,3- dihydrobenzofuran,
18	5-(2-Aminopropyl)-2, 3-dihydrobenzofuran,
19	6-(2-Aminopropyl)-2,3-dihydrobenzofuran,
20	and 7-(2-Aminopropyl)-2,3-dihydrobenzofuran;
21	(12) Methylaminopropylbenzofuran
22	(MAPB): including 4-(2-methylaminopropyl)
23	benzofuran, 5-(2-methylaminopropyl)benzofuran,
24	6-(2-methylaminopropyl)benzofuran
25	and 7-(2-methylaminopropyl)benzofuran.
26	(g) Temporary listing of substances subject to emergency

scheduling. Any material, compound, mixture, or preparation
 that contains any quantity of the following substances:

3 (1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide 4 (benzylfentanyl), its optical isomers, isomers, salts, and 5 salts of isomers;

6 (2) N-[1(2-thienyl) methyl-4-piperidyl]-N-7 phenylpropanamide (thenylfentanyl), its optical isomers, 8 salts, and salts of isomers.

9 (h) Synthetic cathinones. Unless specifically excepted, 10 any chemical compound which is not approved by the United 11 States Food and Drug Administration or, if approved, is not 12 dispensed or possessed in accordance with State or federal including bupropion, structurally derived from 13 law, not 2-aminopropan-1-one by substitution at the 1-position with 14 15 either phenyl, naphthyl, or thiophene ring systems, whether or 16 not the compound is further modified in one or more of the 17 following ways:

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

1 (buphedrone); or

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

8 Any other synthetic cathinone which is not approved by the 9 United States Food and Drug Administration or, if approved, is 10 not dispensed or possessed in accordance with State or federal 11 law.

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

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

20 (Text of Section after amendment by P.A. 103-245)

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

(b) Unless specifically excepted or unless listed in
another schedule, any of the following opiates, including
their isomers, esters, ethers, salts, and salts of isomers,

- 25 - LRB103 35349 RLC 65413 b HB4352 esters, and ethers, whenever the existence of such isomers, 1 2 esters, ethers and salts is possible within the specific chemical designation: 3 4 (1) Acetylmethadol; 5 (1.1) Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-6 7 4-piperidinyl]-N-phenylacetamide); (2) Allylprodine; 8 9 (3) Alphacetylmethadol, except 10 levo-alphacetylmethadol (also known as levo-alpha-11 acetylmethadol, levomethadyl acetate, or LAAM); 12 (4) Alphameprodine; 13 (5) Alphamethadol; (6) Alpha-methylfentanyl 14 (N-(1-alpha-methyl-beta-phenyl) ethyl-4-piperidyl) 15 16 propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-17 propanilido) piperidine; (6.1) Alpha-methylthiofentanyl 18 (N-[1-methyl-2-(2-thienyl)ethyl-19 20 4-piperidinyl]-N-phenylpropanamide); 21 (7) 1-methyl-4-phenyl-4-propionoxypiperidine (MPPP); 22 (7.1) PEPAP 23 (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine); 24 (8) Benzethidine; 25 (9) Betacetylmethadol; 26 (9.1) Beta-hydroxyfentanyl

1	(N-[1-(2-hydroxy-2-phenethyl)-
2	4-piperidinyl]-N-phenylpropanamide);
3	(10) Betameprodine;
4	(11) Betamethadol;
5	(12) Betaprodine;
6	(13) Clonitazene;
7	(14) Dextromoramide;
8	(15) Diampromide;
9	(16) Diethylthiambutene;
10	(17) Difenoxin;
11	(18) Dimenoxadol;
12	(19) Dimepheptanol;
13	(20) Dimethylthiambutene;
14	(21) Dioxaphetylbutyrate;
15	(22) Dipipanone;
16	(23) Ethylmethylthiambutene;
17	(24) Etonitazene;
18	(25) Etoxeridine;
19	(26) Furethidine;
20	(27) Hydroxpethidine;
21	(28) Ketobemidone;
22	(29) Levomoramide;
23	(30) Levophenacylmorphan;
24	(31) 3-Methylfentanyl
25	(N-[3-methyl-1-(2-phenylethyl)-
26	4-piperidyl]-N-phenylpropanamide);

1	(31.1) 3-Methylthiofentanyl
2	(N-[(3-methyl-1-(2-thienyl)ethyl-
3	4-piperidinyl]-N-phenylpropanamide);
4	(32) Morpheridine;
5	(33) Noracymethadol;
6	(34) Norlevorphanol;
7	(35) Normethadone;
8	(36) Norpipanone;
9	(36.1) Para-fluorofentanyl
10	(N-(4-fluorophenyl)-N-[1-(2-phenethyl)-
11	4-piperidinyl]propanamide);
12	(37) Phenadoxone;
13	(38) Phenampromide;
14	(39) Phenomorphan;
15	(40) Phenoperidine;
16	(41) Piritramide;
17	(42) Proheptazine;
18	(43) Properidine;
19	(44) Propiram;
20	(45) Racemoramide;
21	(45.1) Thiofentanyl
22	(N-phenyl-N-[1-(2-thienyl)ethyl-
23	4-piperidinyl]-propanamide);
24	(46) Tilidine;
25	(47) Trimeperidine;
26	(48) Beta-hydroxy-3-methylfentanyl (other name:

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1	N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-
2	N-phenylpropanamide);
3	(49) Furanyl fentanyl (FU-F);
4	(50) Butyryl fentanyl;
5	(51) Valeryl fentanyl;
6	(52) Acetyl fentanyl;
7	(53) Beta-hydroxy-thiofentanyl;
8	(54) 3,4-dichloro-N-[2-
9	(dimethylamino)cyclohexyl]-N-
10	methylbenzamide (U-47700);
11	(55) 4-chloro-N-[1-[2-
12	(4-nitrophenyl)ethyl]-2-piperidinylidene]-
13	benzenesulfonamide (W-18);
14	(56) 4-chloro-N-[1-(2-phenylethyl)
15	-2-piperidinylidene]-benzenesulfonamide (W-15);
16	(57) acrylfentanyl (acryloylfentanyl).
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
21	the 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 lis

(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

HB4352 - 30 - LRB103 35349 RLC 65413 b paragraph only, the term "isomer" includes the optical, 1 2 position and geometric isomers): 3 (1) 3,4-methylenedioxyamphetamine (alpha-methyl, 3, 4-methylenedioxyphenethylamine, 4 5 methylenedioxyamphetamine, MDA); (1.1) Alpha-ethyltryptamine 6 (some trade or other names: etryptamine; 7 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, 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 4 5 plant presently classified botanically as Salvia 6 divinorum, whether growing or not, the seeds thereof, any 7 extract from any part of that plant, and every compound, 8 manufacture, salts, isomers, and salts of isomers whenever 9 the existence of such salts, isomers, and salts of isomers 10 is possible within the specific chemical designation, 11 derivative, mixture, or preparation of that plant, its 12 seeds or 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);

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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:

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

1	<pre>dibenzo (b,d) pyran; Synhexyl;</pre>
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;

- 33 - LRB103 35349 RLC 65413 b HB4352 (26) 2,5-dimethoxy-4-ethylamphetamine 1 2 (another name: DOET); (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine 3 4 (another name: TCPy); 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 (29.1) Benzothiophene analog of phencyclidine. Some 9 10 trade or other names: BTCP or benocyclidine; 11 (29.2) 3-Methoxyphencyclidine (3-MeO-PCP); 12 (30) Bufotenine (some trade or other names: 13 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; 14 5-hydroxy-N,N-dimethyltryptamine; 15 16 N, N-dimethylserotonin; mappine); 17 (31) (Blank); (32) (Blank); 18 19 (33) (Blank); 20 (34) (Blank); 21 (34.5) (Blank); 22 (35) (6aR,10aR) -9- (hydroxymethyl) -6,6-dimethyl-3-23 (2-methyloctan-2-yl)-6a,7, 24 10,10a-tetrahydrobenzo[c]chromen-1-ol 25 Some trade or other names: HU-210; 26 (35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-

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1	dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
2	<pre>tetrahydrobenzo[c]chromen-1-ol, its isomers,</pre>
3	salts, and salts of isomers; Some trade or other
4	names: HU-210, Dexanabinol;
5	(36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-
6	6,6-dimethyl-3-(2-methyloctan-2-yl)-
7	6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol
8	Some trade or other names: HU-211;
9	(37) (Blank);
10	(38) (Blank);
11	(39) (Blank);
12	(40) (Blank);
13	(41) (Blank);
14	(42) Any compound structurally derived from
15	3-(1-naphthoyl)indole or
16	1H-indol-3-yl-(1-naphthyl)methane by substitution at the
17	nitrogen atom of the indole ring by alkyl, haloalkyl,
18	alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide,
19	alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or
20	2-(4-morpholinyl)ethyl whether or not further substituted
21	in the indole ring to any extent, whether or not
22	substituted in the naphthyl ring to any extent. Examples
23	of this structural class include, but are not limited to,
24	JWH-018, AM-2201, JWH-175, JWH-184, and JWH-185;
25	(43) Any compound structurally derived from

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

atom of the pyrrole ring by alkyl, haloalkyl, alkenyl, 1 2 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 3 halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 4 5 in the pyrrole ring to any extent, whether or not substituted in the naphthyl ring to any extent. Examples 6 7 of this structural class include, but are not limited to, JWH-030, JWH-145, JWH-146, JWH-307, and JWH-368; 8

9 (44)Any compound structurally derived from 10 1-(1-naphthylmethyl)indene by substitution at the 11 3-position of the indene ring by alkyl, haloalkyl, 12 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 13 2-(4-morpholinyl)ethyl whether or not further substituted 14 15 in the indene ring to any extent, whether or not 16 substituted in the naphthyl ring to any extent. Examples 17 of this structural class include, but are not limited to, JWH-176; 18

19 (45)compound structurally derived from Any 20 3-phenylacetylindole by substitution at the nitrogen atom indole ring with alkyl, haloalkyl, alkenyl, 21 of the 22 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 23 halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 24 25 in the indole ring to any extent, whether or not 26 substituted in the phenyl ring to any extent. Examples of

1 2 this structural class include, but are not limited to, JWH-167, JWH-250, JWH-251, and RCS-8;

3 (46) Any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at 4 the 5 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, 6 7 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 8 2-(4-morpholinyl)ethyl, whether or not substituted in the 9 cyclohexyl ring to any extent. Examples of this structural 10 class include, but are not limited to, CP 47, 497 and its 11 C8 homologue (cannabicyclohexanol);

12 (46.1)Any compound structurally derived from 13 3-(benzoyl) indole with substitution at the nitrogen atom 14 of the indole ring by an alkyl, haloalkyl, alkenyl, 15 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 16 halide, 1-(N-methyl-2-piperidinyl)methyl, or 17 2-(4-morpholinyl)ethyl group whether or not further substituted in the indole ring to any extent and whether 18 or not substituted in the phenyl ring to any extent. 19 20 Examples of this structural class include, but are not limited to, AM-630, AM-2233, AM-694, Pravadoline (WIN 21 22 48,098), and RCS-4;

(47) (Blank);

23

- 24 (48) (Blank);
- 25 (49) (Blank);
- 26 (50) (Blank);

1	(51) (Blank);
2	(52) (Blank);
3	(53) 2,5-Dimethoxy-4-(n)-propylthio-phenethylamine.
4	Some trade or other names: 2C-T-7;
5	(53.1) 4-ethyl-2,5-dimethoxyphenethylamine. Some
6	<pre>trade or other names: 2C-E;</pre>
7	(53.2) 2,5-dimethoxy-4-methylphenethylamine. Some
8	trade or other names: 2C-D;
9	(53.3) 4-chloro-2,5-dimethoxyphenethylamine. Some
10	<pre>trade or other names: 2C-C;</pre>
11	(53.4) 4-iodo-2,5-dimethoxyphenethylamine. Some trade
12	or other names: 2C-I;
13	(53.5) 4-ethylthio-2,5-dimethoxyphenethylamine. Some
14	trade or other names: 2C-T-2;
15	(53.6) 2,5-dimethoxy-4-isopropylthio-phenethylamine.
16	Some trade or other names: 2C-T-4;
17	(53.7) 2,5-dimethoxyphenethylamine. Some trade or
18	other names: 2C-H;
19	(53.8) 2,5-dimethoxy-4-nitrophenethylamine. Some
20	trade or other names: 2C-N;
21	(53.9) 2,5-dimethoxy-4-(n)-propylphenethylamine. Some
22	trade or other names: 2C-P;
23	(53.10) 2,5-dimethoxy-3,4-dimethylphenethylamine.
24	Some trade or other names: 2C-G;
25	(53.11) The N-(2-methoxybenzyl) derivative of any 2C
26	phenethylamine referred to in subparagraphs (20.1), (53),

(53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7),
 (53.8), (53.9), and (53.10) including, but not limited to,
 25I-NBOMe and 25C-NBOMe;

4

5

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

- (55) (Blank);
- 6 (56) (Blank);
- 7 (57) (Blank);
- 8 (58) (Blank);

9 (59) 3-cyclopropoylindole with substitution at the 10 nitrogen atom of the indole ring by alkyl, haloalkyl, 11 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, 12 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 13 14 on the indole ring to any extent, whether or not 15 substituted on the cyclopropyl ring to any extent: 16 including, but not limited to, XLR11, UR144, FUB-144;

17 (60) 3-adamantoylindole with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, 18 alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, 19 20 alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 21 2-(4-morpholinyl)ethyl, whether or not further substituted 22 on the indole ring to any extent, whether or not substituted on 23 the adamantyl ring to any extent: including, but not limited to, AB-001; 24

25(61)N-(adamantyl)-indole-3-carboxamidewith26substitution at the nitrogen atom of the indole ring by

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haloalkyl, 1 alkyl, alkenyl, cycloalkylmethyl, 2 cycloalkylethyl, aryl halide, alkyl aryl halide, 3 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 4 on the indole ring to any extent, whether or not 5 any extent: 6 substituted on the adamantyl ring to including, but not limited to, APICA/2NE-1, STS-135; 7

N-(adamantyl)-indazole-3-carboxamide 8 (62) with 9 substitution at a nitrogen atom of the indazole ring by 10 alkvl, haloalkyl, alkenyl, cycloalkylmethyl, 11 cycloalkylethyl, aryl halide, alkyl aryl halide, 12 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 13 14 on the indazole ring to any extent, whether or not 15 substituted on the adamantyl ring to any extent: 16 including, but not limited to, AKB48, 5F-AKB48;

17 (63) 1H-indole-3-carboxylic acid 8-quinolinyl ester with substitution at the nitrogen atom of the indole ring 18 19 by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, 20 cycloalkylethyl, aryl halide, alkyl aryl halide, 21 1-(N-methyl-2-piperidinyl)methyl, or 22 2-(4-morpholinyl)ethyl, whether or not further substituted 23 on the indole ring to any extent, whether or not 24 substituted on the quinoline ring to anv extent: 25 including, but not limited to, PB22, 5F-PB22, FUB-PB-22; 26 (64) 3-(1-naphthoyl) indazole with substitution at 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, whether or not substituted on the naphthyl ring to any extent: including, but not limited to, THJ-018, THJ-2201;

8 (65) 2-(1-naphthoyl)benzimidazole with substitution 9 at the nitrogen atom of the benzimidazole ring by alkyl, 10 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 11 aryl halide, alkyl aryl halide, 12 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 13 14 on the benzimidazole ring to any extent, whether or not 15 substituted on the naphthyl ring to any extent: including, 16 but not limited to, FUBIMINA;

17 (66)

18 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indazole-

3-carboxamide with substitution on the nitrogen atom of 19 20 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 on the indazole ring to any extent: including, but not 24 25 limited to, AB-PINACA, AB-FUBINACA, AB-CHMINACA;

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

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

8 (68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-9 indole-3-carboxamide with substitution on the nitrogen 10 atom of the indole ring by alkyl, haloalkyl, alkenyl, 11 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 12 halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted 13 14 on the indole ring to any extent: including, but not 15 limited to, ADBICA, 5F-ADBICA;

16 (69) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indole-17 3-carboxamide with substitution on the nitrogen atom of ring by alkyl, haloalkyl, alkenyl, 18 the indole 19 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 20 halide, 1-(N-methyl-2-piperidinyl)methyl, or 21 2-(4-morpholinyl)ethyl, whether or not further substituted 22 on the indole ring to any extent: including, but not 23 limited 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-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, AMB, 5F-AMB;

2-(1H-indazole-3-carboxamido)-3,3-6 (71)Methvl 7 dimethylbutanoate with substitution on the nitrogen atom 8 indazole ring by alkyl, haloalkyl, alkenyl, of the 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 limited to, 5-fluoro-MDMB-PINACA, MDMB-FUBINACA; 13

14 (72)Methyl 2-(1H-indole-3-carboxamido)-3-15 methylbutanoate with substitution on the nitrogen atom of 16 the indole ring by alkyl, haloalkyl, alkenyl, 17 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, 18 or 2-(4-morpholinyl)ethyl, whether or not further substituted 19 20 on the indazole ring to any extent: including, but not limited to, MMB018, MMB2201, and AMB-CHMICA; 21

(73) Methyl 2-(1H-indole-3-carboxamido)-3,3dimethylbutanoate with substitution on the nitrogen atom
of the indole 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, but not
 limited to, MDMB-CHMICA;

(74)N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-4 5 indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, 6 cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl 7 1-(N-methyl-2-piperidinyl)methyl, 8 halide, or 9 2-(4-morpholinyl)ethyl, whether or not further substituted 10 on the indazole ring to any extent: including, but not 11 limited to, APP-CHMINACA, 5-fluoro-APP-PINACA;

12 (75) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of 13 14 indole ring by alkyl, haloalkyl, alkenvl, the 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 on the indazole ring to any extent: including, but not 18 limited to, APP-PICA and 5-fluoro-APP-PICA; 19

20 (76) 4-Acetoxy-N,N-dimethyltryptamine: trade name 21 4-AcO-DMT;

22 (77) 5-Methoxy-N-methyl-N-isopropyltryptamine: trade 23 name 5-MeO-MIPT;

(78) 4-hydroxy Diethyltryptamine (4-HO-DET);
(79) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);
(80) 4-hydroxy-N, N-diisopropyltryptamine (4-HO-DiPT);

1 (81) 4-hydroxy-N-methyl-N-isopropyltryptamine
2 (4-HO-MiPT);

3

(82) Fluorophenylpiperazine;

4

(83) Methoxetamine;

5 (84) 1-(Ethylamino)-2-phenylpropan-2-one (iso6 ethcathinone).

7 (e) Unless specifically excepted or unless listed in 8 another schedule, any material, compound, mixture, or 9 preparation which contains any quantity of the following 10 substances having a depressant effect on the central nervous 11 system, including its salts, isomers, and salts of isomers 12 whenever the existence of such salts, isomers, and salts of 13 isomers is possible within the specific chemical designation:

14

(1) mecloqualone;

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(2) methaqualone; and

16

(3) gamma hydroxybutyric acid.

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

22

(1) Fenethylline;

23 (2) N-ethylamphetamine;

24 (3) Aminorex (some other names:

25 2-amino-5-phenyl-2-oxazoline; aminoxaphen;

26 4-5-dihydro-5-phenyl-2-oxazolamine) and its

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1	salts, optical isomers, and salts of optical isomers;
2	(4) Methcathinone (some other names:
3	2-methylamino-1-phenylpropan-1-one;
4	Ephedrone; 2-(methylamino)-propiophenone;
5	alpha-(methylamino)propiophenone; N-methylcathinone;
6	methycathinone; Monomethylpropion; UR 1431) and its
7	salts, optical isomers, and salts of optical isomers;
8	(5) Cathinone (some trade or other names:
9	2-aminopropiophenone; alpha-aminopropiophenone;
10	<pre>2-amino-1-phenyl-propanone; norephedrone);</pre>
11	(6) N,N-dimethylamphetamine (also known as:
12	N,N-alpha-trimethyl-benzeneethanamine;
13	N,N-alpha-trimethylphenethylamine);
14	(7) (+ or -) cis-4-methylaminorex ((+ or -) cis-
15	4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);
16	<pre>(8) 3,4-Methylenedioxypyrovalerone (MDPV);</pre>
17	(9) Halogenated amphetamines and
18	methamphetamines - any compound derived from either
19	amphetamine or methamphetamine through the substitution
20	of a halogen on the phenyl ring, including, but not
21	limited to, 2-fluoroamphetamine, 3-
22	fluoroamphetamine and 4-fluoroamphetamine;
23	(10) Aminopropylbenzofuran (APB):
24	including 4-(2-Aminopropyl) benzofuran, 5-
25	(2-Aminopropyl)benzofuran, 6-(2-Aminopropyl)
26	benzofuran, and 7-(2-Aminopropyl) benzofuran;

1	(11) Aminopropyldihydrobenzofuran (APDB):
2	including 4-(2-Aminopropyl)-2,3- dihydrobenzofuran,
3	5-(2-Aminopropyl)-2, 3-dihydrobenzofuran,
4	6-(2-Aminopropyl)-2,3-dihydrobenzofuran,
5	and 7-(2-Aminopropyl)-2,3-dihydrobenzofuran;
6	(12) Methylaminopropylbenzofuran
7	(MAPB): including 4-(2-methylaminopropyl)
8	benzofuran, 5-(2-methylaminopropyl)benzofuran,
9	6-(2-methylaminopropyl)benzofuran
10	and 7-(2-methylaminopropyl)benzofuran.
11	(g) Temporary listing of substances subject to emergency
12	scheduling. Any material, compound, mixture, or preparation
13	that contains any quantity of the following substances:
14	(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
15	(benzylfentanyl), its optical isomers, isomers, salts, and
16	salts of isomers;
17	(2) N-[1(2-thienyl) methyl-4-piperidyl]-N-
18	phenylpropanamide (thenylfentanyl), its optical isomers,
19	salts, and salts of isomers.
20	(h) Synthetic cathinones. Unless specifically excepted,
21	any chemical compound which is not approved by the United
22	States Food and Drug Administration or, if approved, is not
23	dispensed or possessed in accordance with State or federal
24	law, not including bupropion, structurally derived from
25	2-aminopropan-1-one by substitution at the 1-position with
26	either phenyl, naphthyl, or thiophene ring systems, whether or

1 not the compound is further modified in one or more of the 2 following ways:

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

9 (2) by substitution at the 3-position with an acyclic 10 alkyl substituent. Examples of this class include, but are 11 not limited to, 2-methylamino-1-phenylbutan-1-one 12 (buphedrone); or

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

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

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

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1 federal law.

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

8 (1) Benzodiazepine class: A fused 1,4-diazepine and 9 benzene ring structure with a phenyl connected to the 10 1,4-diazepine ring, with any substitution(s) or 11 replacement(s) on the 1,4-diazepine or benzene ring, any 12 substitution(s) on the phenyl ring, or any combination thereof. Examples of this class include but are not 13 14 limited to: Clonazolam, Flualprazolam; or

15 (2) Thienodiazepine class: A fused 1,4-diazepine and 16 thiophene ring structure with a phenyl connected to the 17 with 1,4-diazepine ring, any substitution(s) or replacement(s) on the 1,4-diazepine or thiophene ring, any 18 substitution(s) on the phenyl ring, or any combination 19 thereof. Examples of this class include but are not 20 limited to: Etizolam. 21

22 (Source: P.A. 103-245, eff. 1-1-24.)

(720 ILCS 570/206) (from Ch. 56 1/2, par. 1206)
Sec. 206. (a) The controlled substances listed in this
Section are included in Schedule II.

12

1 (b) Unless specifically excepted or unless listed in 2 another schedule, any of the following substances whether 3 produced directly or indirectly by extraction from substances 4 of vegetable origin, or independently by means of chemical 5 synthesis, or by combination of extraction and chemical 6 synthesis:

7 (1) Opium and opiates, and any salt, compound,
8 derivative or preparation of opium or opiate, excluding
9 apomorphine, dextrorphan, levopropoxyphene, nalbuphine,
10 nalmefene, naloxone, and naltrexone, and their respective
11 salts, but including the following:

- (i) Raw Opium;
- 13 (ii) Opium extracts;
- 14 (iii) Opium fluid extracts;
- 15 (iv) Powdered opium;
- 16 (v) Granulated opium;
- 17 (vi) Tincture of opium;
- 18 (vii) Codeine;
- 19 (viii) Ethylmorphine;
- 20 (ix) Etorphine Hydrochloride;
- 21 (x) Hydrocodone;
- 22 (xi) Hydromorphone;
- 23 (xii) Metopon;
- 24 (xiii) Morphine;
- 25 (xiii.5) 6-Monoacetylmorphine;
- 26 (xiv) Oxycodone;

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- (xv) Oxymorphone;
- 2 (xv.5) Tapentadol;
- 3 (xvi) Thebaine;

(xvii) Thebaine-derived butorphanol.

5 (xviii) Methorphan, except drug products 6 containing dextromethorphan that may be dispensed 7 pursuant to a prescription order of a practitioner and are sold in compliance with the safety and labeling 8 9 standards as set forth by the United States Food and 10 Drug Administration, or drug products containing 11 dextromethorphan that are sold in solid, tablet, 12 liquid, capsule, powder, thin film, or gel form and 13 which are formulated, packaged, and sold in dosages 14 and concentrations for use as an over-the-counter drug 15 product. For the purposes of this Section, 16 "over-the-counter drug product" means a drug that is 17 available to consumers without a prescription and sold in compliance with the safety and labeling standards 18 19 as set forth by the United States Food and Drug 20 Administration.

(2) Any salt, compound, isomer, derivative or
preparation thereof which is chemically equivalent or
identical with any of the substances referred to in
subparagraph (1), but not including the isoquinoline
alkaloids of opium;

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(3) Opium poppy and poppy straw;

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(4) Coca leaves and any salt, compound, isomer, salt 1 2 of an isomer, derivative, or preparation of coca leaves 3 including cocaine or ecgonine, and any salt, compound, isomer, derivative, or preparation thereof which is 4 5 chemically equivalent or identical with any of these substances, but not including decocainized coca leaves or 6 7 extractions of coca leaves which do not contain cocaine or ecgonine (for the purpose of this paragraph, the term 8 9 "isomer" includes optical, positional and geometric 10 isomers);

(5) Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid or powder form which contains the phenanthrine alkaloids of the opium poppy).

(c) Unless specifically excepted or unless listed in another schedule any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, whenever the existence of these isomers, esters, ethers and salts is possible within the specific chemical designation, dextrorphan excepted:

- 20 (1) Alfentanil;
- 21 (1.1) Carfentanil;
- 22 (1.2) Thiafentanyl;
- 23 (2) Alphaprodine;
- 24 (3) Anileridine;
- 25 (4) Bezitramide;

26 (5) Bulk Dextropropoxyphene (non-dosage forms);

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1	(6) Dihydrocodeine;
2	(7) Diphenoxylate;
3	<pre>(8) Fentanyl;</pre>
4	(9) Sufentanil;
5	(9.5) Remifentanil;
6	(10) Isomethadone;
7	(11) (Blank);
8	(12) Levorphanol (Levorphan);
9	(13) Metazocine;
10	(14) Methadone;
11	(15) Methadone-Intermediate,
12	4-cyano-2-dimethylamino-4,4-diphenyl-1-butane;
13	(16) Moramide-Intermediate,
14	2-methyl-3-morpholino-1,1-diphenylpropane-carboxylic
15	acid;
16	(17) Pethidine (meperidine);
17	(18) Pethidine-Intermediate-A,
18	4-cyano-1-methyl-4-phenylpiperidine;
19	(19) Pethidine-Intermediate-B,
20	ethyl-4-phenylpiperidine-4-carboxylate;
21	(20) Pethidine-Intermediate-C,
22	1-methyl-4-phenylpiperidine-4-carboxylic acid;
23	(21) Phenazocine;
24	(22) Piminodine;
25	(23) Racemethorphan;
26	(24) (Blank);

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(25) Levo-alphacetylmethadol (some other names:
 levo-alpha-acetylmethadol, levomethadyl acetate, LAAM).

3 (d) Unless specifically excepted or unless listed in 4 another schedule, any material, compound, mixture, or 5 preparation which contains any quantity of the following 6 substances having a stimulant effect on the central nervous 7 system:

8 (1) Amphetamine, its salts, optical isomers, and salts
9 of its optical isomers;

10 (2) Methamphetamine, its salts, isomers, and salts of
11 its isomers;

(3) Phenmetrazine and its salts;

12

(4) Methylphenidate;

13 14

(5) Lisdexamfetamine.

(e) Unless specifically excepted or unless listed in 15 16 another schedule, any material, compound, mixture, or 17 preparation which contains any quantity of the following substances having a depressant effect on the central nervous 18 system, including its salts, isomers, and salts of isomers 19 20 whenever the existence of such salts, isomers, and salts of 21 isomers is possible within the specific chemical designation:

22

(1) Amobarbital;

23 (2) Secobarbital;

24 (3) Pentobarbital;

25 (4) Pentazocine;

26 (5) Phencyclidine;

- 54 - LRB103 35349 RLC 65413 b HB4352 (6) Gluthethimide; 1 2 (7) (Blank). Unless specifically excepted or unless listed in 3 (f) 4 another schedule, any material, compound, mixture, or 5 preparation which contains any quantity of the following 6 substances: 7 (1)Immediate precursor to amphetamine and 8 methamphetamine: 9 (i) Phenylacetone 10 Some trade or other names: phenyl-2-propanone; 11 P2P; benzyl methyl ketone; methyl benzyl ketone. 12 (2) Immediate precursors to phencyclidine: 13 (i) 1-phenylcyclohexylamine; (ii) 1-piperidinocyclohexanecarbonitrile (PCC). 14 15 (3) Nabilone. 16 (4) Xylazine: N-(2,6-dimethylphenyl)-5,6-dihydro-17 4H-1,3-thiazin-2-amine), including its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, 18 19 whenever the existence of such isomers, esters, ethers, 20 and salts is possible within the specific chemical 21 designation as a Schedule II controlled substance. 22 (g) Clonazolam. 23 (Source: P.A. 100-368, eff. 1-1-18.)

24 Section 95. No acceleration or delay. Where this Act makes 25 changes in a statute that is represented in this Act by text that is not yet or no longer in effect (for example, a Section represented by multiple versions), the use of that text does not accelerate or delay the taking effect of (i) the changes made by this Act or (ii) provisions derived from any other Public Act.