

HOUSE BILL No. 1196

DIGEST OF INTRODUCED BILL

Citations Affected: IC 35-41-1-26.3; IC 35-48-2-4.

Synopsis: Synthetic drugs (including "bath salts"). Adds additional chemical compounds (including some compounds sold as "bath salts") to the definition of synthetic cannabinoids, and expands the definition of synthetic cannabinoids to encompass certain chemical compounds that are structurally related to synthetic cannabinoids.

Effective: July 1, 2012.

Smith M, Yarde

January 9, 2012, read first time and referred to Committee on Courts and Criminal Code.

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Second Regular Session 117th General Assembly (2012)

PRINTING CODE. Amendments: Whenever an existing statute (or a section of the Indiana Constitution) is being amended, the text of the existing provision will appear in this style type, additions will appear in **this style type**, and deletions will appear in ~~this style type~~.

Additions: Whenever a new statutory provision is being enacted (or a new constitutional provision adopted), the text of the new provision will appear in **this style type**. Also, the word **NEW** will appear in that style type in the introductory clause of each SECTION that adds a new provision to the Indiana Code or the Indiana Constitution.

Conflict reconciliation: Text in a statute in *this style type* or ~~this style type~~ reconciles conflicts between statutes enacted by the 2011 Regular Session of the General Assembly.

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HOUSE BILL No. 1196

A BILL FOR AN ACT to amend the Indiana Code concerning criminal law and procedure.

Be it enacted by the General Assembly of the State of Indiana:

1 SECTION 1. IC 35-41-1-26.3, AS ADDED BY P.L.182-2011,
2 SECTION 10, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
3 JULY 1, 2012]: Sec. 26.3. "Synthetic cannabinoid" means:
4 (1) a substance containing one (1) or more of the following
5 chemical compounds, **including an analogue or homologue of**
6 **the compound:**
7 (1) (A) JWH-015 ((2-Methyl-1-propyl-1H-
8 indol-3-yl)-1-naphthalenylmethanone).
9 (2) (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
10 (3) (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
11 (4) (D) JWH-073
12 (naphthalen-1-yl-(1-butylyndol-3-yl)methanone).
13 (5) (E) JWH-081 (4-methoxynaphthalen- 1-yl-
14 (1-pentylindol- 3-yl)methanone).
15 (6) (F) JWH-122
16 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
17 (7) ~~JWH-200 (1-(2-morpholin-4-ylethyl)indol-3-yl)-~~



- 1 naphthalen-1-ylmethanone):
 2 **(G) JWH-200**
 3 **((1-(2-morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-ylm**
 4 **ethanone).**
 5 **(8) (H) JWH-250**
 6 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
 7 **(9) (I) JWH-251** (1-pentyl-3-(2-methylphenylacetyl)indole).
 8 **(10) (J) JWH-398**
 9 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
 10 **(11) (K) HU-210** ((6aR,10aR)- 9-(Hydroxymethyl)-
 11 6,6-dimethyl- 3-(2-methyloctan-2-yl)-
 12 6a,7,10,10a-tetrahydrobenzo [c]chromen- 1-ol).
 13 **(12) (L) HU-211** ((6aS,10aS)-9-(Hydroxymethyl)-
 14 6,6-dimethyl- 3-(2-methyloctan-2-yl)-
 15 6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
 16 **(13) (M) HU-308** ([(1R,2R,5R)-2-[2,6-dimethoxy-4-
 17 (2-methyloctan- 2-yl)phenyl]-
 18 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
 19 **(14) HU-331** ((3-hydroxy-2- [(1R,6R)-3-methyl-6-
 20 (1-methylethenyl)-2 -cyclohexen-1-yl]-5
 21 -pentyl-2,5-cyclohexadiene-1,4-dione).
 22 **(N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-**
 23 **(1-methylethenyl)-2 -cyclohexen-1-yl]-5**
 24 **-pentyl-2,5-cyclohexadiene-1,4-dione).**
 25 **(15) (O) CP 55,940**
 26 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]-
 27 5- (2-methyloctan-2-yl)phenol).
 28 **(16) (P) CP 47,497** (2-[(1R,3S)-3-hydroxycyclohexyl]- 5-
 29 (2-methyloctan-2-yl)phenol) and its homologues, **or**
 30 **2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)**
 31 **phenol), where side chain n=5, and homologues where**
 32 **side chain n-4,6, or 7.**
 33 **(17) (Q) WIN 55212-2**
 34 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
 35 pyrrolo [1,2,3-de)- 1,4- benzoxazin-
 36 6-yl]-1-naphthalenylmethanone).
 37 **(18) (R) RCS-4** ((4-methoxyphenyl)
 38 (1-pentyl-1H-indol-3-yl)methanone).
 39 **(19) (S) RCS-8** (1-(1-(2-cyclohexylethyl)-1H-
 40 indol-3-yl)-2-(2-methoxyphenyl)ethanone).
 41 **(20) (T) 4-Methylmethcathinone.** Other name: mephedrone.
 42 **(21) (U) 3,4-Methylenedioxy-methcathinone.** Other name:

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- 1 methylone.
 2 (22) (V) Fluoromethcathinone.
 3 (23) (W) 4-Methoxymethcathinone. Other name:
 4 methedrone.
 5 (24) (X) 4-Ethylmethcathinone (4-EMC).
 6 (25) (Y) Methylendioxyprovalerone. Other name: MDPV.
 7 (Z) JWH-007, or
 8 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
 9 (AA) JWH-098, or
 10 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
 11 (BB) JWH-164, or
 12 1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
 13 (CC) JWH-210, or
 14 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
 15 (DD) JWH-201, or
 16 1-pentyl-3-(4-methoxyphenylacetyl)indole.
 17 (EE) JWH-203, or
 18 1-pentyl-3-(2-chlorophenylacetyl)indole.
 19 (FF) AM-694, or
 20 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
 21 (GG) CP 50,556-1, or
 22 [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenyl
 23 pentan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthri
 24 din-1-yl] acetate.
 25 (HH) Dimethylheptylpyran, or DMHP.
 26 (II) 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP.
 27 (2) Any compound structurally derived from
 28 3-(1-naphthoyl)indole or
 29 1H-indol-3-yl-(1-naphthyl)methane by substitution at the
 30 nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl,
 31 cycloalkylmethyl, cycloalkylethyl,
 32 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
 33 group, whether or not further substituted in the indole ring
 34 to any extent and whether or not substituted in the
 35 naphthyl ring to any extent.
 36 (3) Any compound structurally derived from
 37 3-(1-naphthoyl)pyrrole by substitution at the nitrogen
 38 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl,
 39 cycloalkylmethyl, cycloalkylethyl,
 40 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
 41 group, whether or not further substituted in the pyrrole
 42 ring to any extent and whether or not substituted in the

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- 1 naphthyl ring to any extent.
2 (4) Any compound structurally derived from
3 1-(1-naphthylmethyl)indene by substitution at the
4 3-position of the indene ring by alkyl, haloalkyl, alkenyl,
5 cycloalkylmethyl, cycloalkylethyl,
6 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
7 group, whether or not further substituted in the indene ring
8 to any extent and whether or not substituted in the
9 naphthyl ring to any extent.
10 (5) Any compound structurally derived from
11 3-phenylacetylindole by substitution at the nitrogen atom of
12 the indole ring with alkyl, haloalkyl, alkenyl,
13 cycloalkylmethyl, cycloalkylethyl,
14 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
15 group, whether or not further substituted in the indole ring
16 to any extent and whether or not substituted in the phenyl
17 ring to any extent.
18 (6) Any compound structurally derived from
19 2-(3-hydroxycyclohexyl)phenol by substitution at the
20 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl,
21 cycloalkylmethyl, cycloalkylethyl,
22 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
23 group, whether or not substituted in the cyclohexyl ring to
24 any extent.
25 (7) Any compound containing a 3-(benzoyl)indole structure
26 with substitution at the nitrogen atom of the indole ring by
27 alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
28 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl
29 group, whether or not further substituted in the indole ring
30 to any extent and whether or not substituted in the phenyl
31 ring to any extent.
32 (8) Any compound, except bupropion or a compound listed
33 under a different schedule, structurally derived from
34 2-aminopropan-1-one by substitution at the 1-position with
35 either phenyl, naphthyl, or thiophene ring systems, whether
36 or not the compound is further modified:
37 (A) by substitution in the ring system to any extent with
38 alkyl, alkylenedioxy, alkoxy, haloalkyl, hydroxyl, or
39 halide substituents, whether or not further substituted in
40 the ring system by one or more other univalent
41 substituents;
42 (B) by substitution at the 3-position with an acyclic alkyl

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1 **substituent;**
 2 **(C) by substitution at the 2-amino nitrogen atom with**
 3 **alkyl, dialkyl, benzyl, or methoxybenzyl groups; or**
 4 **(D) by inclusion of the 2-amino nitrogen atom in a cyclic**
 5 **structure.**

6 SECTION 2. IC 35-48-2-4, AS AMENDED BY P.L.182-2011,
 7 SECTION 12, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 8 JULY 1, 2012]: Sec. 4. (a) The controlled substances listed in this
 9 section are included in schedule I.

10 (b) Opiates. Any of the following opiates, including their isomers,
 11 esters, ethers, salts, and salts of isomers, esters, and ethers, unless
 12 specifically excepted by rule of the board or unless listed in another
 13 schedule, whenever the existence of these isomers, esters, ethers, and
 14 salts is possible within the specific chemical designation:

15 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-
 16 piperidinyl]-N-phenylacetamide) (9815)

17 Acetylmethadol (9601)

18 Allylprodine (9602)

19 Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
 20 thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)

21 Alphacetylmethadol (9603)

22 Alphameprodine (9604)

23 Alphamethadol (9605)

24 Alphamethylfentanyl (9814)

25 Benzethidine (9606)

26 Beta-hydroxy-3-methylfentanyl (9831). Other name:

27 N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl
 28]-N-phenylpropanamide

29 Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-
 30 phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)

31 Betacetylmethadol (9607)

32 Betameprodine (9608)

33 Betamethadol (9609)

34 Betaprodine (9611)

35 Clonitazene (9612)

36 Dextromoramide (9613)

37 Diampromide (9615)

38 Diethylthiambutene (9616)

39 Difenoxin (9168)

40 Dimenoxadol (9617)

41 Dimepheptanol (9618)

42 Dimethylthiambutene (9619)

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1	Dioxaphetyl butyrate (9621)
2	Dipipanone (9622)
3	Ethylmethylthiambutene (9623)
4	Etonitazene (9624)
5	Etoxidine (9625)
6	Furethidine (9626)
7	Hydroxypethidine (9627)
8	Ketobemidone (9628)
9	Levomoramide (9629)
10	Levophenacymorphan (9631)
11	3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-
12	piperidyl]-N-phenyl-propanamide](9813)
13	3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
14	piperidinyl]-N-phenylpropanamide) (9833)
15	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
16	Morpheridine (9632)
17	N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
18	(benzylfentanyl), including any isomers, salts, or salts of
19	isomers (9818)
20	N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide
21	(thenylfentanyl), including any isomers, salts, or salts of
22	isomers (9834)
23	Noracymethadol (9633)
24	Norlevorphanol (9634)
25	Normethadone (9635)
26	Norpipanone (9636)
27	Para-fluorofentanyl (N-(4-fluorophenyl)-N-
28	[1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
29	Phenadoxone (9637)
30	Phenampromide (9638)
31	Phenomorphane (9647)
32	Phenoperidine (9641)
33	PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
34	Piritramide (9642)
35	Proheptazine (9643)
36	Propiridine (9644)
37	Propiram (9649)
38	Racemoramide (9645)
39	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
40	piperidinyl]-propanamide) (9835)
41	Tilidine (9750)
42	Trimeperidine (9646)

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1 (c) Opium derivatives. Any of the following opium derivatives, their
 2 salts, isomers, and salts of isomers, unless specifically excepted by rule
 3 of the board or unless listed in another schedule, whenever the
 4 existence of these salts, isomers, and salts of isomers is possible within
 5 the specific chemical designation:

6 Acetorphine (9319)
 7 Acetyldihydrocodeine (9051)
 8 Benzylmorphine (9052)
 9 Codeine methylbromide (9070)
 10 Codeine-N-Oxide (9053)
 11 Cyprenorphine (9054)
 12 Desomorphine (9055)
 13 Dihydromorphine (9145)
 14 Drotebanol (9335)
 15 Etorphine (except hydrochloride salt) (9056)
 16 Heroin (9200)
 17 Hydromorphinol (9301)
 18 Methyldesorphine (9302)
 19 Methyl dihydromorphine (9304)
 20 Morphine methylbromide (9305)
 21 Morphine methylsulfonate (9306)
 22 Morphine-N-Oxide (9307)
 23 Myrophine (9308)
 24 Nicocodeine (9309)
 25 Nicomorphine (9312)
 26 Normorphine (9313)
 27 Pholcodine (9314)
 28 Thebacon (9315)

29 (d) Hallucinogenic substances. Any material, compound, mixture,
 30 or preparation which contains any quantity of the following
 31 hallucinogenic, psychedelic, or psychogenic substances, their salts,
 32 isomers, and salts of isomers, unless specifically excepted by rule of
 33 the board or unless listed in another schedule, whenever the existence
 34 of these salts, isomers, and salts of isomers is possible within the
 35 specific chemical designation:

36 (1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name:
 37 TCPy.
 38 (2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade
 39 or other names: 4-Bromo-2,
 40 5-Dimethoxy-a-methylphenethylamine; 4-Bromo-2, 5-DMA.
 41 (3) 4-Bromo-2, 5-dimethoxyphenethylamine (7392). Some trade
 42 or other names:

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- 1 2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;
 2 alpha-desmethyl DOB; 2C-B, Nexus.
 3 (4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name:
 4 DOET.
 5 (5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348).
 6 Other name: 2C-T-7.
 7 (6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other
 8 names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.
 9 (7) 4-Methoxyamphetamine (7411). Some trade or other
 10 names: 4-Methoxy-a-methylphenethylamine;
 11 Paramethoxyamphetamine; PMA.
 12 (8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401).
 13 Other Name: MMDA.
 14 (9) 5-Methoxy-N, N-diisopropyltryptamine, including any
 15 isomers, salts, or salts of isomers (7439). Other name:
 16 5-MeO-DIPT.
 17 (10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade
 18 and other names: 4-methyl-2,
 19 5-dimethoxy-a-methylphenethylamine; DOM; and STP.
 20 (11) 3, 4-methylenedioxy amphetamine (7400). Other name:
 21 MDA.
 22 (12) 3,4-methylenedioxy-N-ethylamphetamine (7404). Other
 23 names: N-ethyl-alpha-methyl-3,4(methylenedioxy)
 24 phenethylamine; N-ethyl MDA; MDE; and MDEA.
 25 (13) 3, 4-methylenedioxymethamphetamine (MDMA) (7405).
 26 (14) 3, 4, 5-trimethoxy amphetamine (7390). Other name:
 27 TMA.
 28 (15) Alpha-ethyltryptamine (7249). Some trade and other
 29 names: Etryptamine; Monase;
 30 [alpha]-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)
 31 indole; [alpha]-ET; and AET.
 32 (16) Alpha-methyltryptamine (7432). Other name: AMT.
 33 (17) Bufotenine (7433). Some trade and other names:
 34 3-(B-Dimethylaminoethyl)-5-hydroxyindole;
 35 3-(2-dimethylaminonethyl)-5-indolol; N, N-dimethylserotonin;
 36 5-hydroxy-N, N-dimethyltryptamine; mappine.
 37 (18) Diethyltryptamine (7434). Some trade or other names: N,
 38 N-Diethyltryptamine; DET.
 39 (19) Dimethyltryptamine (7435). Some trade or other names:
 40 DMT.
 41 (20) Ibogaine (7260). Some trade and other names: 7-Ethyl-6,
 42 6b, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6,

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- 1 9-methano-5H-pyrido (1', 2': 1, 2, azepino 4, 5-b) indole;
 2 tabernanthe iboga.
 3 (21) Lysergic acid diethylamide (7315). Other name: LSD.
 4 (22) Marijuana (7360).
 5 (23) Mescaline (7381).
 6 (24) Parahexyl (7374). Some trade or other names:
 7 3-Hexyl-1-hydroxy-7, 8, 9, 10-Tetrahydro-6, 6,
 8 9-trimethyl-6H-dibenzo (b,d) pyran; Snyhexyl.
 9 (25) Peyote (7415), including:
 10 (A) all parts of the plant that are classified botanically as
 11 lophophora williamsii lemaire, whether growing or not;
 12 (B) the seeds thereof;
 13 (C) any extract from any part of the plant; and
 14 (D) every compound, manufacture, salt, derivative, mixture,
 15 or preparation of the plant, its seeds, or extracts.
 16 (26) N-ethyl-3-piperidyl benzilate (7482). Other name: DMZ.
 17 (27) N-hydroxy-3,4-methylenedioxyamphetamine (7402).
 18 Other names: N-hydroxy-alpha-methyl-3,4
 19 (methylenedioxy)phenethylamine; and N-hydroxy MDA.
 20 (28) N-methyl-3-piperidyl benzilate (7484). Other name: LBJ.
 21 (29) Psilocybin (7437).
 22 (30) Psilocyn (7438).
 23 (31) Tetrahydrocannabinols (7370), including synthetic
 24 equivalents of the substances contained in the plant, or in the
 25 resinous extractives of Cannabis, sp. and synthetic substances,
 26 derivatives, and their isomers with similar chemical structure
 27 and pharmacological activity such as:
 28 (A) π^1 cis or trans tetrahydrocannabinol, and their optical
 29 isomers;
 30 (B) π^6 cis or trans tetrahydrocannabinol, and their optical
 31 isomers; and
 32 (C) π^3_4 cis or trans tetrahydrocannabinol, and their optical
 33 isomers.
 34 Since nomenclature of these substances is not internationally
 35 standardized, compounds of these structures, regardless of
 36 numerical designation of atomic positions are covered. Other
 37 name: THC.
 38 (32) Ethylamine analog of phencyclidine (7455). Some trade or
 39 other names: N-Ethyl-1-phenylcyclohexylamine;
 40 (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl)
 41 ethylamine; cyclohexamine; PCE.
 42 (33) Pyrrolidine analog of phencyclidine (7458). Some trade or

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- 1 other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCP,_y; PHP.
 2 (34) Thiophene analog of phencyclidine (7470). Some trade or
 3 other names: 1-(1-(2-thienyl) cyclohexyl) piperidine; 2-Thienyl
 4 Analog of Phencyclidine; TPCP.
 5 (35) Synthetic cannabinoids (as defined in IC 35-41-1-26.3).
 6 including a substance containing one (1) or more of the
 7 following chemical compounds:
 8 (A) JWH-015 ((2-Methyl-1-propyl-1H-
 9 indol-3-yl)-1-naphthalenylmethanone).
 10 (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
 11 (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
 12 (D) JWH-073 (naphthalen-1-yl-
 13 (1-butylindol-3-yl)methanone).
 14 (E) JWH-081 (4-methoxynaphthalen-1-yl- (1-pentylindol-
 15 3-yl)methanone).
 16 (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
 17 (G) JWH-200 (1-(2-morpholin-4-ylethyl)indol-3-yl)-
 18 naphthalen-1-ylmethanone).
 19 (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
 20 (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
 21 (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
 22 (K) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-
 23 6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
 24 tetrahydrobenzo [c]chromen-1-ol).
 25 (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-
 26 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo
 27 [c]chromen-1-ol).
 28 (M) HU-308 ((1R,2R,5R)-2-[2,6-dimethoxy-4-
 29 (2-methyloctan-2-yl)phenyl]-
 30 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
 31 (N) HU-331 ((3-hydroxy-2-[(1R,6R)-3-methyl-6-
 32 (1-methylethenyl)-2-cyclohexen-1-yl]-5-
 33 -pentyl-2,5-cyclohexadiene-1,4-dione).
 34 (O) CP 55,940 (2-[(1R,2R,5R)-5-hydroxy-
 35 2-(3-hydroxypropyl) cyclohexyl]-5-
 36 (2-methyloctan-2-yl)phenol).
 37 (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]-5-
 38 (2-methyloctan-2-yl)phenol) and its homologues.
 39 (Q) WIN 55212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-
 40 (4-morpholinylmethyl) pyrrolo [1,2,3-de]-1,4- benzoxazin-
 41 6-yl]-1-naphthalenylmethanone).
 42 (R) RCS-4 ((4-methoxyphenyl)

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- 1 (1-pentyl-1H-indol-3-yl)methanone):
 2 (S) RCS-8 (1-(1-(2-cyclohexylethyl)-
 3 1H-indol-3-yl)-2-(2-methoxyphenyl)ethanone):
 4 (T) 4-Methylmethcathinone. Other name: mephedrone.
 5 (U) 3,4-Methylenedioxymethcathinone. Other name:
 6 methylone.
 7 (V) Fluoromethcathinone.
 8 (W) 4-Methoxymethcathinone. Other name: methedrone.
 9 (X) 4-Ethylmethcathinone. Other name: 4-EMC.
 10 (Y) Methylenedioxyprovalerone. Other name: MDPV.
 11 (36) Salvia divinorum or salvinorin A, including:
 12 (A) all parts of the plant that are classified botanically as salvia
 13 divinorum, whether growing or not;
 14 (B) the seeds of the plant;
 15 (C) any extract from any part of the plant; and
 16 (D) every compound, manufacture, salt, derivative, mixture, or
 17 preparation of the plant, its seeds, or extracts.
 18 (e) Depressants. Unless specifically excepted in a rule adopted by
 19 the board or unless listed in another schedule, any material, compound,
 20 mixture, or preparation which contains any quantity of the following
 21 substances having a depressant effect on the central nervous system,
 22 including its salts, isomers, and salts of isomers whenever the existence
 23 of such salts, isomers, and salts of isomers is possible within the
 24 specific chemical designation:
 25 Gamma-hydroxybutyric acid (other names include GHB;
 26 gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium
 27 oxybate; sodium oxybutyrate) (2010)
 28 Mecloqualone (2572)
 29 Methaqualone (2565)
 30 (f) Stimulants. Unless specifically excepted or unless listed in
 31 another schedule, any material, compound, mixture, or preparation that
 32 contains any quantity of the following substances having a stimulant
 33 effect on the central nervous system, including its salts, isomers, and
 34 salts of isomers:
 35 ([+/-] cis-4-methylaminorex (([+/-]cis-4,5-
 36 dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)
 37 Aminorex (1585). Other names: aminoxaphen;
 38 2-amino-5-phenyl-2-oxazoline; or
 39 4,5-dihydro-5-phenyl-2-oxazolamine.
 40 Cathinone (1235). Some trade or other names:
 41 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone;
 42 2-aminopropiophenone; and norephedrone.

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- 1 Fenethylamine (1503).
- 2 N-Benzylpiperazine (7493). Other names: BZP; and
- 3 1-benzylpiperazine.
- 4 N-ethylamphetamine (1475)
- 5 Methcathinone (1237) Some other trade names:
- 6 2-Methylamino-1-Phenylpropan-1-one; Ephedrone;
- 7 Monomethylpropion; UR 1431.
- 8 N, N-dimethylamphetamine (1480). Other names: N,
- 9 N-alpha-trimethyl-benzeneethanamine; and N,
- 10 N-alpha-trimethylphenethylamine.

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