



March 15, 2017

ENGROSSED HOUSE BILL No. 1019

DIGEST OF HB 1019 (Updated March 14, 2017 9:46 am - DI 84)

Citations Affected: IC 35-31.5; IC 35-48.

Synopsis: Controlled substances. Adds the substance U-47700 to Schedule I. Adds Etizolam to Schedule I. Adds the chemical description of buphedrone.

Effective: July 1, 2017.

Ellington, Negele, Klinker, Hatfield

(SENATE SPONSORS — MERRITT, HEAD, GROOMS)

January 4, 2017, read first time and referred to Committee on Courts and Criminal Code.
February 16, 2017, amended, reported — Do Pass.
February 20, 2017, read second time, ordered engrossed. Engrossed.
February 21, 2017, read third time, passed. Yeas 90, nays 0.

SENATE ACTION

February 23, 2017, read first time and referred to Committee on Corrections and Criminal Law.
March 14, 2017, reported favorably — Do Pass.

EH 1019—LS 6164/DI 106



March 15, 2017

First Regular Session of the 120th General Assembly (2017)

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 2016 Regular Session of the General Assembly.

ENGROSSED HOUSE BILL No. 1019

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 35-31.5-2-321, AS AMENDED BY P.L.168-2016,
2 SECTION 10, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
3 JULY 1, 2017]: Sec. 321. "Synthetic drug" means:
4 (1) a substance containing one (1) or more of the following
5 chemical compounds, including an analog of the compound:
6 (A) JWH-015 ((2-Methyl-1-propyl-1H-
7 indol-3-yl)-1-naphthalenylmethanone).
8 (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
9 (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
10 (D) JWH-073
11 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).
12 (E) JWH-081 (4-methoxynaphthalen- 1-yl- (1-pentylindol-
13 3-yl)methanone).
14 (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
15 (G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-
16 naphthalen-1-yl-methanone).
17 (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).

EH 1019—LS 6164/DI 106



- 1 (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
 2 (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
 3 (K) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-
 4 3-(2-methyloctan-2-yl)-
 5 6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
 6 (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-
 7 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo
 8 [c]chromen-1-ol).
 9 (M) HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-
 10 (2-methyloctan-2-yl)phenyl]-
 11 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
 12 (N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-
 13 (1-methylethenyl)-2-cyclohexen-1-yl]-5
 14 -pentyl-2,5-cyclohexadiene-1,4-dione).
 15 (O) CP 55,940
 16 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-
 17 5-(2-methyloctan-2-yl)phenol).
 18 (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]-5-
 19 (2-methyloctan-2-yl)phenol) and its homologues, or
 20 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)
 21 phenol, where side chain n=5, and homologues where side
 22 chain n=4, 6, or 7.
 23 (Q) WIN 55212-2
 24 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
 25 pyrrolo [1,2,3-de)-1,4-benzoxazin-
 26 6-yl]-1-naphthalenylmethanone).
 27 (R) RCS-4 ((4-methoxyphenyl)
 28 (1-pentyl-1H-indol-3-yl)methanone).
 29 (S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-
 30 indol-3-yl)-2-(2-methoxyphenyl)ethanone).
 31 (T) 4-Methylmethcathinone. Other name: mephedrone.
 32 (U) 3,4-Methylenedioxy-methcathinone. Other name:
 33 methylone.
 34 (V) Fluoromethcathinone.
 35 (W) 4-Methoxymethcathinone. Other name: methedrone.
 36 (X) 4-Ethylmethcathinone (4-EMC).
 37 (Y) Methylenedioxypropylone. Other name: MDPV.
 38 (Z) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
 39 (AA) JWH-098, or
 40 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
 41 (BB) JWH-164, or
 42 1-pentyl-3-(7-methoxy-1-naphthoyl)indole.



1	(CC) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
2	(DD) JWH-201, or
3	1-pentyl-3-(4-methoxyphenylacetyl)indole.
4	(EE) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole.
5	(FF) AM-694, or
6	1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
7	(GG) CP 50,556-1, or
8	[(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpe
9	ntan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1
10	-yl] acetate.
11	(HH) Dimethylheptylpyran, or DMHP.
12	(II) 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP.
13	(JJ) 6-APB [6-(2-aminopropyl)benzofuran].
14	(LL) 7-hydroxymitragynine.
15	(MM) α -PPP [α -pyrrolidinopropiophenone].
16	(NN) α -PVP (desmethylpyrovalerone).
17	(OO) AM-251.
18	(PP) AM-1241.
19	(QQ) AM-2201.
20	(RR) AM-2233.
21	(SS) Buphedrone (α-methylamino-butyrophenone (MABP)).
22	(TT) Butylone.
23	(UU) CP-47,497-C7.
24	(VV) CP-47,497-C8.
25	(WW) Desoxypipradol.
26	(XX) Ethylone.
27	(YY) Eutylone.
28	(ZZ) Flephedrone.
29	(AAA) JWH-011.
30	(BBB) JWH-020.
31	(CCC) JWH-022.
32	(DDD) JWH-030.
33	(EEE) JWH-182.
34	(FFF) JWH-302.
35	(GGG) MDAI [5,6-methylenedioxy-2-aminoindane].
36	(HHH) Mitragynine.
37	(III) Naphyrone.
38	(JJJ) Pentedrone.
39	(LLL) Pentylone.
40	(MMM) Methoxetamine
41	[2-(3-methoxyphenyl)-2-(ethylamino)- cyclohexanone].
42	(NNN) A796,260 [1-(2-morpholin-4-ylethyl)-1H-indol-3-yl]-



1 (2,2,3,3-tetramethylcyclopropyl)methanone].
 2 (OOO) AB-001[(1s,3s)-adamantan-1-yl]
 3 (1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
 4 (1-adamantoyl)indole].
 5 (PPP) AM-356 [Methanandamide].
 6 (QQQ) AM 1248 [1-[(1-methyl-2- piperidinyl) methyl]-
 7 1H-indol-3-yl] tricyclo[3.3.1.1^{3,7}] dec-1-yl-methanone]or
 8 [(1-[(N-methylpiperindin-2-yl)
 9 Methyl]-3-(Adamant-1-oyl)indole)].
 10 (RRR) AM 2233 Azepane isomer [(2-iodophenyl)
 11 (1-(1-methylazepan-3-yl)- 1H-indol-3-yl)methanone].
 12 (SSS) CB-13 [1-Naphthalenyl
 13 [4-(pentyoxy)- 1-naphthalenyl]methanone].
 14 (TTT) UR-144 [(1-pentyl-1H-indol-3-yl)
 15 (2,2,3,3-tetramethylcyclopropyl)-methanone].
 16 (UUU) URB 597 [(3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl)-
 17 cyclohexylcarbamate].
 18 (VVV) URB602 [[1,1'-biphenyl]- 3-yl-carbamic acid,
 19 cyclohexyl ester].
 20 (WWW) URB 754 [6-methyl-2-[(4-methylphenyl)
 21 amino]-1-benzoxazin-4-one].
 22 (XXX) XLR-11 or 5-fluoro UR-144
 23 (1-(5-fluoropentyl)-1H-indol-3-yl)
 24 (2,2,3,3-tetramethylcyclopropyl)methanone].
 25 (YYY) AKB48 (Other names include:
 26 N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide;
 27 1-pentyl-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-indazole-3-
 28 carboxamide).
 29 (ZZZ) 25I-NBOMe (Other names include:
 30 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-
 31 benzeneethanamine);
 32 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
 33 methyl]ethanamine).
 34 (AAAA) 2C-C-NBOMe (Other names include: 25C-NBOMe;
 35 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
 36 methyl]ethanamine;
 37 2,5-Dimethoxy-4-chloro-N-(2-methoxybenzyl)
 38 phenethylamine).
 39 (BBBB) 2NE-1 (Other names include: 1-Pentyl-3-
 40 (1-adamantylamido)indole).
 41 (CCCC) STS-135 (Other names include:
 42 N-Adamantyl-1-fluoropentylindole-3- carboxamide



1 (1-5-fluoropentyl)-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-
2 indole-3-carboxamide).
3 (DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinyl
4 ester-1H-indole-2-carboxylic acid).
5 (EEEE) 5-Fluoro-PB-22 (Other names include:
6 1-(5-Fluoropentyl)-8-quinolinyl ester-1H-indole-3-carboxylic
7 acid).
8 (FFFF) Benocyclidine (Other names include: BCP, BTCP, and
9 Benzothiophenylcyclohexylpiperidine).
10 (GGGG) 25B-NBOMe (Other names include: 2C-B-NBOMe
11 and 4-Bromo-2,
12 5-dimethoxy-N-[(2-Methoxyphenyl)methyl]
13 benzeneethanamine).
14 (HHHH) APB (Other names include; (2-Aminopropyl)
15 Benzofuran).
16 (III) AB-PINACA
17 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
18 indazole-3-carboxamide).
19 (JJJJ) AB-FUBINACA
20 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-
21 1H-indazole-3-carboxamide).
22 (KKKK) ADB-PINACA
23 (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-in
24 daole-3-carboxamide).
25 (LLLL) Fluoro ADBICA (N-(1-Amino-3,3-
26 dimethyl-1-oxobutan-2-yl)-(fluoropentyl)-1H-indole-3-
27 carboxamide).
28 (MMMM) APDB (Other names include: -EMA,
29 -Desoxy-MDA, and (2-Aminopropyl)-2,3-
30 dihydrobenzofuran).
31 (NNNN) THJ-2201 (Other names include: AM2201 indazole
32 analog, Fluoropentyl-JWH-018 indazole, and
33 5-Fluoro-THJ-018).
34 (OOOO) AM 2201 benzimidazole analog (Other names
35 include: FUBIMINA, FTHJ, and (1-(5-fluoropentyl)-1H-
36 benzo[d]imidazol-2-yl)(naphthalene-1-yl)methanone).
37 (PPPP) MN-25 (Other names include: 7-methoxy-1-
38 [2-(4-morpholinyl)ethyl]-N-[1S, 2S, 4R)-1,3,3-
39 trimethylbicyclo[2.2.1]hept-2-yl]-1H-indole-3-carboxamide
40 and UR-12).
41 (QQQQ) FUB-PB-22 (Other names include:
42 Quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).



- 1 (RRRR) FUD-PB-22 (Other names include:
 2 Naphthalen-1-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 3 (SSSS) 5-Fluoro-AB-PINACA (Other names include:
 4 AB-PINACA 5-fluoro analog and N-(1-amino-3-methyl-
 5 oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
 6 carboxamide).
 7 (TTTT) 4-MePPP (Other names include:
 8 4-methyl-alpha-pyrrolidinopropiophenone).
 9 (UUUU) alpha-PBP (Other names include:
 10 Alpha-pyrrolidinobutiophenone).
 11 (VVVV) AB-CHMINACA (Other names include:
 12 (N-[1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethy
 13 l)-1H-indazole-3-carboxamide).
 14 (WWWW) Acetyl fentanyl (Other names include:
 15 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide).
 16 (2) Any compound structurally derived from
 17 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by
 18 substitution at the nitrogen atom of the indole ring by alkyl,
 19 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 20 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 21 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 22 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 23 or not further substituted in the indole ring to any extent and
 24 whether or not substituted in the naphthyl ring to any extent.
 25 (3) Any compound structurally derived from 3-(1-naphthoyl)
 26 pyrrole by substitution at the nitrogen atom of the pyrrole ring by
 27 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
 28 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,
 29 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl,
 30 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
 31 group, whether or not further substituted in the pyrrole ring to any
 32 extent and whether or not substituted in the naphthyl ring to any
 33 extent.
 34 (4) Any compound structurally derived from
 35 1-(1-naphthylmethyl)indene by substitution at the 3-position of
 36 the indene ring by alkyl, haloalkyl, cyanoalkyl, alkenyl,
 37 c y c l o a l k y l m e t h y l , c y c l o a l k y l e t h y l ,
 38 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 39 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 40 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 41 or not further substituted in the indene ring to any extent and
 42 whether or not substituted in the naphthyl ring to any extent.



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



- 1 cycloalkylmethyl, cycloalkylethyl,
2 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl) ethyl,
3 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
4 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
5 or not further substituted in the indole ring to any extent and
6 whether or not substituted in the tetramethylcyclopropyl ring to
7 any extent.
- 8 (10) Any compound containing a N-(1-adamantyl)-
9 1H-indazole-3-carboxamide structure with substitution at the
10 nitrogen atom of the indazole ring by an alkyl, haloalkyl,
11 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
13 1 - (N - m e t h y l - 2 - p y r r o l i d i n y l) m e t h y l ,
14 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
15 group, whether or not further substituted at the nitrogen atom of
16 the carboxamide to any extent, whether or not further substituted
17 in the indazole ring to any extent, and whether or not further
18 substituted on the adamantyl ring system to any extent. An
19 example of this structural class includes AKB48.
- 20 (11) Any compound containing a N-(1-adamantyl)-
21 1H-indole-3-carboxamide structure with substitution at the
22 nitrogen atom of the indole ring by an alkyl, haloalkyl,
23 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
24 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
25 1 - (N - m e t h y l - 2 - p y r r o l i d i n y l) m e t h y l ,
26 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
27 group, whether or not further substituted at the nitrogen atom of
28 the carboxamide to any extent, whether or not further substituted
29 in the indole ring to any extent, and whether or not further
30 substituted on the adamantyl ring system to any extent. An
31 example of this structural class includes STS-135.
- 32 (12) Any compound containing a 3-(1-adamantoyl)indole
33 structure with substitution at the nitrogen atom of the indole ring
34 by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
35 cycloalkylethyl, 1-(N-methyl-2- piperidinyl)methyl, or
36 2-(4-morpholinyl)ethyl, 1-(N-methyl-2- pyrrolidinyl)methyl,
37 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
38 group, whether or not further substituted on the adamantyl ring
39 system to any extent. An example of this structural class includes
40 AM-1248.
- 41 (13) Any compound determined to be a synthetic drug by rule
42 adopted under IC 25-26-13-4.1.



1 SECTION 2. IC 35-48-2-4, AS AMENDED BY P.L.168-2016,
 2 SECTION 11, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 3 JULY 1, 2017]: Sec. 4. (a) The controlled substances listed in this
 4 section are included in schedule I.

5 (b) Opiates. Any of the following opiates, including their isomers,
 6 esters, ethers, salts, and salts of isomers, esters, and ethers, unless
 7 specifically excepted by rule of the board or unless listed in another
 8 schedule, whenever the existence of these isomers, esters, ethers, and
 9 salts is possible within the specific chemical designation:

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

12 Acetylmethadol (9601)

13 Allylprodine (9602)

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

16 Alphacetylmethadol (9603)

17 Alphameprodine (9604)

18 Alphamethadol (9605)

19 Alphamethylfentanyl (9814)

20 Benzethidine (9606)

21 Beta-hydroxy-3-methylfentanyl (9831). Other name:
 22 N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl
 23]-N-phenylpropanamide

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

26 Betacetylmethadol (9607)

27 Betameprodine (9608)

28 Betamethadol (9609)

29 Betaprodine (9611)

30 Clonitazene (9612)

31 Dextromoramide (9613)

32 Diampromide (9615)

33 Diethylthiambutene (9616)

34 Difenoxin (9168)

35 Dimenoxadol (9617)

36 Dimepheptanol (9618)

37 Dimethylthiambutene (9619)

38 Dioxaphetyl butyrate (9621)

39 Dipipanone (9622)

40 Ethylmethylthiambutene (9623)

41 Etonitazene (9624)

42 Etoxidine (9625)



- 1 Furethidine (9626)
 2 Hydroxypethidine (9627)
 3 Ketobemidone (9628)
 4 Levomoramide (9629)
 5 Levophenacymorphan (9631)
 6 3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-
 7 piperidyl]-N-phenyl-propanamide](9813)
 8 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
 9 piperidinyl]-N-phenylpropanamide) (9833)
 10 MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
 11 Morpheridine (9632)
 12 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl),
 13 including any isomers, salts, or salts of isomers (9818)
 14 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide
 15 (thenylfentanyl), including any isomers, salts, or salts of isomers
 16 (9834)
 17 Noracymethadol (9633)
 18 Norlevorphanol (9634)
 19 Normethadone (9635)
 20 Norpitanone (9636)
 21 Para-fluorofentanyl (N-(4-fluorophenyl)-N-
 22 [1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
 23 Phenadoxone (9637)
 24 Phenampromide (9638)
 25 Phenomorphan (9647)
 26 Phenoperidine (9641)
 27 PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
 28 Piritramide (9642)
 29 Proheptazine (9643)
 30 Properidine (9644)
 31 Propiram (9649)
 32 Racemoramide (9645)
 33 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
 34 piperidinyl]-propanamide) (9835)
 35 Tilidine (9750)
 36 Trimeperidine (9646)
 37 **U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-**
 38 **N-methyl- benzamide)**
 39 (c) Opium derivatives. Any of the following opium derivatives, their
 40 salts, isomers, and salts of isomers, unless specifically excepted by rule
 41 of the board or unless listed in another schedule, whenever the
 42 existence of these salts, isomers, and salts of isomers is possible within



- 1 the specific chemical designation:
- 2 Acetorphine (9319)
- 3 Acetyldihydrocodeine (9051)
- 4 Benzylmorphine (9052)
- 5 Codeine methylbromide (9070)
- 6 Codeine-N-Oxide (9053)
- 7 Cyprenorphine (9054)
- 8 Desomorphine (9055)
- 9 Dihydromorphine (9145)
- 10 Drotebanol (9335)
- 11 Etorphine (except hydrochloride salt) (9056)
- 12 Heroin (9200)
- 13 Hydromorphanol (9301)
- 14 Methyldesorphine (9302)
- 15 Methylhydromorphine (9304)
- 16 Morphine methylbromide (9305)
- 17 Morphine methylsulfonate (9306)
- 18 Morphine-N-Oxide (9307)
- 19 Myrophine (9308)
- 20 Nicocodeine (9309)
- 21 Nicomorphine (9312)
- 22 Normorphine (9313)
- 23 Pholcodine (9314)
- 24 Thebacon (9315)
- 25 (d) Hallucinogenic substances. Unless specifically excepted or
- 26 unless listed in another schedule, any material, compound, mixture, or
- 27 preparation which contains any quantity of the following
- 28 hallucinogenic, psychedelic, or psychogenic substances, their salts,
- 29 isomers, and salts of isomers whenever the existence of these salts,
- 30 isomers, and salts of isomers is possible within the specific chemical
- 31 designation (for purposes of this subsection only, the term "isomer"
- 32 includes the optical, position, and geometric isomers):
- 33 (1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name:
- 34 TCPy.
- 35 (2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or
- 36 other names: 4-Bromo-2, 5-Dimethoxy- α -methylphenethylamine;
- 37 4-Bromo-2, 5-DMA.
- 38 (3) 4-Bromo-2, 5-dimethoxyphenethylamine (7392). Some trade
- 39 or other names:
- 40 2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;
- 41 α -desmethyl DOB; 2C-B, Nexus.
- 42 (4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name:



- 1 DOET.
- 2 (5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348).
- 3 Other name: 2C-T-7.
- 4 (6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other
- 5 names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.
- 6 (7) 4-Methoxyamphetamine (7411). Some trade or other names:
- 7 4-Methoxy-a-methylphenethylamine; Paramethoxyamphetamine;
- 8 PMA.
- 9 (8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401). Other
- 10 Name: MMDA.
- 11 (9) 5-Methoxy-N, N-diisopropyltryptamine, including any
- 12 isomers, salts, or salts of isomers (7439). Other name:
- 13 5-MeO-DIPT.
- 14 (10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade
- 15 and other names: 4-methyl-2,
- 16 5-dimethoxy-a-methylphenethylamine; DOM; and STP.
- 17 (11) 3, 4-methylenedioxy amphetamine (7400). Other name:
- 18 MDA.
- 19 (12) 3,4-methylenedioxy-N-ethylamphetamine (7404). Other
- 20 names: N-ethyl-alpha-methyl-3,4(methylenedioxy)
- 21 phenethylamine; N-ethyl MDA; MDE; and MDEA.
- 22 (13) 3, 4-methylenedioxymethamphetamine (MDMA) (7405).
- 23 (14) 3, 4, 5-trimethoxy amphetamine (7390). Other name: TMA.
- 24 (15) Alpha-ethyltryptamine (7249). Some trade and other names:
- 25 Etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine;
- 26 3-(2-aminobutyl) indole; [alpha]-ET; and AET.
- 27 (16) Alpha-methyltryptamine (7432). Other name: AMT.
- 28 (17) Bufotenine (7433). Some trade and other names:
- 29 3-(B-Dimethylaminoethyl)-5-hydroxyindole;
- 30 3-(2-dimethylaminonethyl)-5-indolol; N, N-dimethylserotonin;
- 31 5-hydroxy-N, N-dimethyltryptamine; mappine.
- 32 (18) Diethyltryptamine (7434). Some trade or other names: N,
- 33 N-Diethyltryptamine; DET.
- 34 (19) Dimethyltryptamine (7435). Some trade or other names:
- 35 DMT.
- 36 (20) Ibogaine (7260). Some trade and other names: 7-Ethyl-6, 6b,
- 37 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6, 9-methano-5H-pyrido
- 38 (1', 2': 1, 2, azepino 4, 5-b) indole; tabernanthe iboga.
- 39 (21) Lysergic acid diethylamide (7315). Other name: LSD.
- 40 (22) Marijuana (7360).
- 41 (23) Mescaline (7381).
- 42 (24) Parahexyl (7374). Some trade or other names:



- 1 3-Hexyl-1-hydroxy-7, 8, 9, 10-Tetrahydro-6, 6,
 2 9-trimethyl-6H-dibenzo (b,d) pyran; Snyhexyl.
 3 (25) Peyote (7415), including:
 4 (A) all parts of the plant that are classified botanically as
 5 lophophora williamsii lemaire, whether growing or not;
 6 (B) the seeds thereof;
 7 (C) any extract from any part of the plant; and
 8 (D) every compound, manufacture, salt, derivative, mixture, or
 9 preparation of the plant, its seeds, or extracts.
 10 (26) N-ethyl-3-piperidyl benzilate (7482). Other name: DMZ.
 11 (27) N-hydroxy-3,4-methylenedioxyamphetamine (7402). Other
 12 names: N-hydroxy-alpha-methyl-3,4
 13 (methylenedioxy)phenethylamine; and N-hydroxy MDA.
 14 (28) N-methyl-3-piperidyl benzilate (7484). Other name: LBJ.
 15 (29) Psilocybin (7437).
 16 (30) Psilocyn (7438).
 17 (31) Tetrahydrocannabinols (7370), including synthetic
 18 equivalents of the substances contained in the plant, or in the
 19 resinous extractives of Cannabis, sp. and synthetic substances,
 20 derivatives, and their isomers with similar chemical structure and
 21 pharmacological activity such as:
 22 (A) π^1 cis or trans tetrahydrocannabinol, and their optical
 23 isomers;
 24 (B) π^6 cis or trans tetrahydrocannabinol, and their optical
 25 isomers; and
 26 (C) π^3_4 cis or trans tetrahydrocannabinol, and their optical
 27 isomers.
 28 Since nomenclature of these substances is not internationally
 29 standardized, compounds of these structures, regardless of
 30 numerical designation of atomic positions are covered. Other
 31 name: THC.
 32 (32) Ethylamine analog of phencyclidine (7455). Some trade or
 33 other names: N-Ethyl-1-phenylcyclohexylamine;
 34 (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl)
 35 ethylamine; cyclohexamine; PCE.
 36 (33) Pyrrolidine analog of phencyclidine (7458). Some trade or
 37 other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCP,_y; PHP.
 38 (34) Thiophene analog of phencyclidine (7470). Some trade or
 39 other names: 1-(1-(2-thienyl) cyclohexyl) piperidine; 2-Thienyl
 40 Analog of Phencyclidine; TPCP.
 41 (35) Salvia divinorum or salvinorin A, including:
 42 (A) all parts of the plant that are classified botanically as salvia



- 1 divinorum, whether growing or not;
 2 (B) the seeds of the plant;
 3 (C) any extract from any part of the plant; and
 4 (D) every compound, manufacture, salt, derivative, mixture, or
 5 preparation of the plant, its seeds, or extracts.
- 6 (36) 5-Methoxy-N,N-Dimethyltryptamine. Some trade or other
 7 names: 5-methoxy-3-[2- (dimethylamino)ethyl]indole;
 8 5-MeO-DMT.
- 9 (37) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E).
 10 (38) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D).
 11 (39) 2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C).
 12 (40) 2-(4-Iodo-2,5-dimethoxyphenyl) ethanamine (2C-I).
 13 (41) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl] ethanamine (2C-T-2).
 14 (42) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl] ethanamine
 15 (2C-T-4).
 16 (43) 2-(2,5-Dimethoxyphenyl) ethanamine (2C-H).
 17 (44) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N).
 18 (45) 2-(2,5-Dimethoxy-4-(n)-propylphenyl) ethanamine (2C-P).
- 19 (e) Depressants. Unless specifically excepted in a rule adopted by
 20 the board or unless listed in another schedule, any material, compound,
 21 mixture, or preparation which contains any quantity of the following
 22 substances having a depressant effect on the central nervous system,
 23 including its salts, isomers, and salts of isomers whenever the existence
 24 of such salts, isomers, and salts of isomers is possible within the
 25 specific chemical designation:
- 26 **Etizolam (4-(2- chlorophenyl)-2- ethyl-9- methyl- 6H-**
 27 **thieno[3,2-f] [1,2,4] triazolo[4,3-a] [1,4diazepine) (other names**
 28 **include: Etilaam, Etizest, Depas, Etizola, Sedekopan, and**
 29 **Pasaden)**
- 30 Gamma-hydroxybutyric acid (other names include GHB;
 31 gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium
 32 oxybate; sodium oxybutyrate) (2010)
 33 Mecloqualone (2572)
 34 Methaqualone (2565)
- 35 (f) Stimulants. Unless specifically excepted or unless listed in
 36 another schedule, any material, compound, mixture, or preparation that
 37 contains any quantity of the following substances having a stimulant
 38 effect on the central nervous system, including its salts, isomers, and
 39 salts of isomers:
- 40 ([+/-]) cis-4-methylaminorex (([+/-])cis-4,5-
 41 dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)
 42 Aminorex (1585). Other names: aminoxaphen;



- 1 2-amino-5-phenyl-2-oxazoline; or
2 4,5-dihydro-5-phenyl-2-oxazolamine.
3 Cathinone (1235). Some trade or other names:
4 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone;
5 2-aminopropiophenone; and norephedrone.
6 Fenethylamine (1503).
7 N-Benzylpiperazine (7493). Other names: BZP; and
8 1-benzylpiperazine.
9 N-ethylamphetamine (1475).
10 Methcathinone (1237) Some other trade names:
11 2-Methylamino-1-Phenylpropan-1-one; Ephedrone;
12 Monomethylpropion; UR 1431.
13 N, N-dimethylamphetamine (1480). Other names: N,
14 N-alpha-trimethyl-benzeneethanamine; and N,
15 N-alpha-trimethylphenethylamine.
16 (g) Synthetic drugs as defined in IC 35-31.5-2-321.



COMMITTEE REPORT

Mr. Speaker: Your Committee on Courts and Criminal Code, to which was referred House Bill 1019, has had the same under consideration and begs leave to report the same back to the House with the recommendation that said bill be amended as follows:

Page 3, line 21, delete "." and insert "**(α -methylamino-butyrophenone (MABP)).**".

Page 6, delete line 16.

Page 9, after line 1, begin a new paragraph and insert:

"SECTION 1. IC 35-48-2-4, AS AMENDED BY P.L.168-2016, SECTION 11, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2017]: Sec. 4. (a) The controlled substances listed in this section are included in schedule I.

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

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

Acetylmethadol (9601)

Allylprodine (9602)

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

Alphacetylmethadol (9603)

Alphameprodine (9604)

Alphamethadol (9605)

Alphamethylfentanyl (9814)

Benzethidine (9606)

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

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

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

Betacetylmethadol (9607)

Betameprodine (9608)

Betamethadol (9609)

Betaprodine (9611)

Clonitazene (9612)

Dextromoramide (9613)

Diampromide (9615)



Diethylthiambutene (9616)
Difenoxin (9168)
Dimenoxadol (9617)
Dimepheptanol (9618)
Dimethylthiambutene (9619)
Dioxaphetyl butyrate (9621)
Dipipanone (9622)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxidine (9625)
Furethidine (9626)
Hydroxypethidine (9627)
Ketobemidone (9628)
Levomoramide (9629)
Levophenacymorphan (9631)
3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide](9813)
3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9833)
MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
Morpheridine (9632)
N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), including any isomers, salts, or salts of isomers (9818)
N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thienylfentanyl), including any isomers, salts, or salts of isomers (9834)
Noracymethadol (9633)
Norlevorphanol (9634)
Normethadone (9635)
Norpipanone (9636)
Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
Phenadoxone (9637)
Phenampramide (9638)
Phenomorphane (9647)
Phenoperidine (9641)
PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
Piritramide (9642)
Proheptazine (9643)
Properidine (9644)
Propiram (9649)
Racemoramide (9645)



Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide) (9835)

Tilidine (9750)

Trimeperidine (9646)

U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-N-methyl- benzamide)

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

Acetorphine (9319)

Acetyldihydrocodeine (9051)

Benzylmorphine (9052)

Codeine methylbromide (9070)

Codeine-N-Oxide (9053)

Cyprenorphine (9054)

Desomorphine (9055)

Dihydromorphine (9145)

Drotebanol (9335)

Etorphine (except hydrochloride salt) (9056)

Heroin (9200)

Hydromorphinol (9301)

Methyl-desorphine (9302)

Methyldihydromorphine (9304)

Morphine methylbromide (9305)

Morphine methylsulfonate (9306)

Morphine-N-Oxide (9307)

Myrophine (9308)

Nicocodeine (9309)

Nicomorphine (9312)

Normorphine (9313)

Pholcodine (9314)

Thebacon (9315)

(d) Hallucinogenic substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following hallucinogenic, psychedelic, or psychogenic substances, their salts, isomers, and salts of isomers whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation (for purposes of this subsection only, the term "isomer" includes the optical, position, and geometric isomers):

EH 1019—LS 6164/DI 106



- (1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name: TCPy.
- (2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or other names: 4-Bromo-2, 5-Dimethoxy-a-methylphenethylamine; 4-Bromo-2, 5-DMA.
- (3) 4-Bromo-2, 5-dimethoxyphenethylamine (7392). Some trade or other names:
2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;
alpha-desmethyl DOB; 2C-B, Nexus.
- (4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name: DOET.
- (5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348). Other name: 2C-T-7.
- (6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.
- (7) 4-Methoxyamphetamine (7411). Some trade or other names: 4-Methoxy-a-methylphenethylamine; Paramethoxyamphetamine; PMA.
- (8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401). Other Name: MMDA.
- (9) 5-Methoxy-N, N-diisopropyltryptamine, including any isomers, salts, or salts of isomers (7439). Other name: 5-MeO-DIPT.
- (10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade and other names: 4-methyl-2, 5-dimethoxy-a-methylphenethylamine; DOM; and STP.
- (11) 3, 4-methylenedioxy amphetamine (7400). Other name: MDA.
- (12) 3,4-methylenedioxy-N-ethylamphetamine (7404). Other names: N-ethyl-alpha-methyl-3,4(methylenedioxy) phenethylamine; N-ethyl MDA; MDE; and MDEA.
- (13) 3, 4-methylenedioxymethamphetamine (MDMA) (7405).
- (14) 3, 4, 5-trimethoxy amphetamine (7390). Other name: TMA.
- (15) Alpha-ethyltryptamine (7249). Some trade and other names: Etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; [alpha]-ET; and AET.
- (16) Alpha-methyltryptamine (7432). Other name: AMT.
- (17) Bufotenine (7433). Some trade and other names: 3-(B-Dimethylaminoethyl)-5-hydroxyindole;
3-(2-dimethylaminonethyl)-5-indolol; N, N-dimethylserotonin;
5-hydroxy-N, N-dimethyltryptamine; mappine.
- (18) Diethyltryptamine (7434). Some trade or other names: N,



N-Diethyltryptamine; DET.

(19) Dimethyltryptamine (7435). Some trade or other names: DMT.

(20) Ibogaine (7260). Some trade and other names: 7-Ethyl-6, 6b, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6, 9-methano-5H-pyrido (1', 2': 1, 2, azepino 4, 5-b) indole; tabernanthe iboga.

(21) Lysergic acid diethylamide (7315). Other name: LSD.

(22) Marijuana (7360).

(23) Mescaline (7381).

(24) Parahexyl (7374). Some trade or other names: 3-Hexyl-1-hydroxy-7, 8, 9, 10-Tetrahydro-6, 6, 9-trimethyl-6H-dibenzo (b,d) pyran; Snyhexyl.

(25) Peyote (7415), including:

(A) all parts of the plant that are classified botanically as *lophophora williamsii* lemaire, whether growing or not;

(B) the seeds thereof;

(C) any extract from any part of the plant; and

(D) every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or extracts.

(26) N-ethyl-3-piperidyl benzilate (7482). Other name: DMZ.

(27) N-hydroxy-3,4-methylenedioxyamphetamine (7402). Other names: N-hydroxy-alpha-methyl-3,4

(methylenedioxy)phenethylamine; and N-hydroxy MDA.

(28) N-methyl-3-piperidyl benzilate (7484). Other name: LBJ.

(29) Psilocybin (7437).

(30) Psilocyn (7438).

(31) Tetrahydrocannabinols (7370), including synthetic equivalents of the substances contained in the plant, or in the resinous extractives of *Cannabis*, sp. and synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity such as:

(A) π^1 cis or trans tetrahydrocannabinol, and their optical isomers;

(B) π^6 cis or trans tetrahydrocannabinol, and their optical isomers; and

(C) π^3_4 cis or trans tetrahydrocannabinol, and their optical isomers.

Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions are covered. Other name: THC.

(32) Ethylamine analog of phencyclidine (7455). Some trade or



other names: N-Ethyl-1-phenylcyclohexylamine; (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl) ethylamine; cyclohexamine; PCE.

(33) Pyrrolidine analog of phencyclidine (7458). Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCP_y; PHP.

(34) Thiophene analog of phencyclidine (7470). Some trade or other names: 1-(1-(2-thienyl) cyclohexyl) piperidine; 2-Thienyl Analog of Phencyclidine; TPCP.

(35) *Salvia divinorum* or salvinorin A, including:

(A) all parts of the plant that are classified botanically as *salvia divinorum*, whether growing or not;

(B) the seeds of the plant;

(C) any extract from any part of the plant; and

(D) every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or extracts.

(36) 5-Methoxy-N,N-Dimethyltryptamine. Some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT.

(37) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E).

(38) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D).

(39) 2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C).

(40) 2-(4-Iodo-2,5-dimethoxyphenyl) ethanamine (2C-I).

(41) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl] ethanamine (2C-T-2).

(42) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl] ethanamine (2C-T-4).

(43) 2-(2,5-Dimethoxyphenyl) ethanamine (2C-H).

(44) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N).

(45) 2-(2,5-Dimethoxy-4-(n)-propylphenyl) ethanamine (2C-P).

(e) Depressants. Unless specifically excepted in a rule adopted by the board or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including 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:

Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f][1,2,4]triazolo[4,3-a][1,4diazepine) (other names include: Etilaam, Etizest, Depas, Etizola, Sedekopan, and Pasaden)

Gamma-hydroxybutyric acid (other names include GHB; gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate) (2010)



Mecloqualone (2572)

Methaqualone (2565)

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

([+/-]) cis-4-methylaminorex (([+/-])cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)

Aminorex (1585). Other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine.

Cathinone (1235). Some trade or other names: 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone; 2-aminopropiophenone; and norephedrone.

Fenethylamine (1503).

N-Benzylpiperazine (7493). Other names: BZP; and 1-benzylpiperazine.

N-ethylamphetamine (1475).

Methcathinone (1237) Some other trade names: 2-Methylamino-1-Phenylpropan-1-one; Ephedrone; Monomethylpropion; UR 1431.

N, N-dimethylamphetamine (1480). Other names: N, N-alpha-trimethyl-benzeneethanamine; and N, N-alpha-trimethylphenethylamine.

(g) Synthetic drugs as defined in IC 35-31.5-2-321."

Renumber all SECTIONS consecutively.

and when so amended that said bill do pass.

(Reference is to HB 1019 as introduced.)

WASHBURN

Committee Vote: yeas 8, nays 0.



COMMITTEE REPORT

Madam President: The Senate Committee on Corrections and Criminal Law, to which was referred House Bill No. 1019, has had the same under consideration and begs leave to report the same back to the Senate with the recommendation that said bill DO PASS.

(Reference is to HB1019 as printed February 17, 2017.)

YOUNG M, Chairperson

Committee Vote: Yeas 7, Nays 0

