



March 29, 2019

ENGROSSED SENATE BILL No. 631

DIGEST OF SB 631 (Updated March 27, 2019 2:24 pm - DI 131)

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

Synopsis: Drug classifications and drug schedules. Adds numerous substances to the definition of "synthetic drug". Adds epidiolex and brivaracetam to schedule V. Specifies that dronabinol is a schedule II controlled substance only in oral solution. Defines "fentanyl related substance" and adds it to schedule I. Moves certain fentanyl related substances from the definition of "synthetic drugs" in schedule II to schedule I. Adds Thiafentanil to schedule II. Makes conforming changes.

Effective: July 1, 2019.

Young M, Freeman, Bohacek

(HOUSE SPONSOR — MCNAMARA)

January 15, 2019, read first time and referred to Committee on Rules and Legislative Procedure.

January 16, 2019, amended; reassigned to Committee on Corrections and Criminal Law.

February 7, 2019, amended, reported favorably — Do Pass.

February 11, 2019, read second time, ordered engrossed. Engrossed.

February 12, 2019, read third time, passed. Yeas 47, nays 1.

HOUSE ACTION

March 7, 2019, read first time and referred to Committee on Courts and Criminal Code.

March 28, 2019, amended, reported — Do Pass.

ES 631—LS 7594/DI 13



March 29, 2019

First Regular Session of the 121st General Assembly (2019)

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 2018 Regular and Special Session of the General Assembly.

ENGROSSED SENATE BILL No. 631

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 35-31.5-2-130.5 IS ADDED TO THE INDIANA
2 CODE AS A **NEW** SECTION TO READ AS FOLLOWS
3 [EFFECTIVE JULY 1, 2019]: **Sec. 130.5. "Fentanyl related**
4 **substance", for purposes of IC 35-48, has the meaning set forth in**
5 **IC 35-48-1-16.6.**

6 SECTION 2. IC 35-31.5-2-321, AS AMENDED BY P.L.89-2018,
7 SECTION 1, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
8 JULY 1, 2019]: Sec. 321. "Synthetic drug" means:

9 (1) a substance containing one (1) or more of the following
10 chemical compounds, including an analog of the compound:

- 11 (A) JWH-015 ((2-Methyl-1-propyl-1H-
12 indol-3-yl)-1-naphthalenylmethanone).
13 (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
14 (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
15 (D) JWH-073
16 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).
17 (E) JWH-081 (4-methoxynaphthalen- 1-yl- (1-pentylindol-

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- 1 3-yl)methanone).
 2 (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
 3 (G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-
 4 naphthalen-1-yl-methanone).
 5 (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
 6 (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
 7 (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
 8 (K) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-
 9 3-(2-methyloctan-2-yl)-
 10 6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
 11 (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-
 12 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo
 13 [c]chromen-1-ol).
 14 (M) HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-
 15 (2-methyloctan-2-yl)phenyl]-
 16 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl]methanol).
 17 (N) HU-331 (3-hydroxy-2-[(1R,6R)-3-methyl-6-
 18 (1-methylethenyl)-2-cyclohexen-1-yl]-5-
 19 pentyl-2,5-cyclohexadiene-1,4-dione).
 20 (O) CP 55,940
 21 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-
 22 5-(2-methyloctan-2-yl)phenol).
 23 (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]-5-
 24 (2-methyloctan-2-yl)phenol) and its homologues, or
 25 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)
 26 phenol, where side chain n=5, and homologues where side
 27 chain n=4, 6, or 7.
 28 (Q) WIN 55212-2
 29 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
 30 pyrrolo [1,2,3-de]-1,4-benzoxazin-
 31 6-yl]-1-naphthalenylmethanone).
 32 (R) RCS-4 ((4-methoxyphenyl)
 33 (1-pentyl-1H-indol-3-yl)methanone).
 34 (S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-
 35 indol-3-yl)-2-(2-methoxyphenyl)ethanone).
 36 (T) 4-Methylmethcathinone. Other name: mephedrone.
 37 (U) 3,4-Methylenedioxymethcathinone. Other name:
 38 methylone.
 39 (V) Fluoromethcathinone.
 40 (W) 4-Methoxymethcathinone. Other name: methedrone.
 41 (X) 4-Ethylmethcathinone (4-EMC).
 42 (Y) Methylenedioxyprovalerone. Other name: MDPV.



1	(Z) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
2	(AA) JWH-098, or
3	1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
4	(BB) JWH-164, or
5	1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
6	(CC) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
7	(DD) JWH-201, or
8	1-pentyl-3-(4-methoxyphenylacetyl)indole.
9	(EE) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole.
10	(FF) AM-694, or
11	1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
12	(GG) CP 50,556-1, or
13	[(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpe
14	ntan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1
15	-yl] acetate.
16	(HH) Dimethylheptylpyran, or DMHP.
17	(II) 4-Methyl- α -pyrrolidinobutiophenone, or MPBP.
18	(JJ) 6-APB [6-(2-aminopropyl)benzofuran].
19	(LL) 7-hydroxymitragynine.
20	(MM) α -PPP [α -pyrrolidinopropiophenone].
21	(NN) α -PVP (desmethylpyrovalerone).
22	(OO) AM-251.
23	(PP) AM-1241.
24	(QQ) AM-2201.
25	(RR) AM-2233.
26	(SS) Buphedrone (α -methylamino-butyrophenone (MABP)).
27	(TT) Butylone.
28	(UU) CP-47,497-C7.
29	(VV) CP-47,497-C8.
30	(WW) Desoxypipradol.
31	(XX) Ethylone.
32	(YY) Eutylone.
33	(ZZ) Flephedrone.
34	(AAA) JWH-011.
35	(BBB) JWH-020.
36	(CCC) JWH-022.
37	(DDD) JWH-030.
38	(EEE) JWH-182.
39	(FFF) JWH-302.
40	(GGG) MDAI [5,6-methylenedioxy-2-aminoindane].
41	(HHH) Mitragynine.
42	(III) Naphyrone.



1	(JJJ) Pentedrone.
2	(LLL) Pentylone.
3	(MMM) Methoxetamine
4	[2-(3-methoxyphenyl)-2-(ethylamino)- cyclohexanone].
5	(NNN) A796,260
6	[1-(2-morpholin-4-ylethyl)-1H-indol-3-yl]-
7	(2,2,3,3-tetramethylcyclopropyl)methanone].
8	(OOO) AB-001[(1s,3s)-adamantan-1-yl]
9	(1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
10	(1-adamantoyl)indole].
11	(PPP) AM-356 [Methanandamide].
12	(QQQ) AM 1248 [1-[(1-methyl-2- piperidinyl) methyl]-
13	1H-indol-3-yl] tricyclo[3.3.1.1 ^{3,7}] dec-1-yl-methanone]or
14	[(1-[(N-methylpiperindin-2-yl)
15	Methyl]-3-(Adamant-1-oyl)indole].
16	(RRR) AM 2233 Azepane isomer [(2-iodophenyl)
17	(1-(1-methylazepan-3-yl)- 1H-indol-3-yl)methanone].
18	(SSS) CB-13 [1-Naphthalenyl
19	[4-(pentyoxy)- 1-naphthalenyl]methanone].
20	(TTT) UR-144 [(1-pentyl-1H-indol-3-yl)
21	(2,2,3,3-tetramethylcyclopropyl)-methanone].
22	(UUU) URB 597 [(3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl)-
23	cyclohexylcarbamate].
24	(VVV) URB602 [[1,1'-biphenyl]- 3-yl-carbamic acid,
25	cyclohexyl ester].
26	(WWW) URB 754 [6-methyl-2-[(4-methylphenyl)
27	amino]-1-benzoxazin-4-one].
28	(XXX) XLR-11 or 5-fluoro UR-144
29	(1-(5-fluoropentyl)-1H-indol-3-yl)
30	(2,2,3,3-tetramethylcyclopropyl)methanone].
31	(YYY) AKB48 (Other names include:
32	N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide;
33	1-pentyl-N-tricyclo[3.3.1.1 ^{3,7}]dec-1-yl-1H-indazole-3-
34	carboxamide).
35	(ZZZ) 25I-NBOMe (Other names include:
36	4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-
37	benzeneethanamine);
38	2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
39	methyl]ethanamine).
40	(AAAA) 2C-C-NBOMe (Other names include:
41	25C-NBOMe;
42	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)



1 methyl]ethanamine;
 2 2,5-Dimethoxy-4-chloro-N-(2-methoxybenzyl)
 3 phenethylamine).
 4 (BBBB) 2NE-1 (Other names include: 1-Pentyl-3-
 5 (1-adamantylamido)indole).
 6 (CCCC) STS-135 (Other names include:
 7 N-Adamantyl-1-fluoropentylindole-3- carboxamide
 8 (1-5-fluoropentyl)-N-tricyclo[3.3.1.1.3.7]dec-1-yl-1H-
 9 indole-3-carboxamide).
 10 (DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinyl
 11 ester-1H-indole-2-carboxylic acid).
 12 (EEEE) 5-Fluoro-PB-22 (Other names include:
 13 1-(5-Fluoropentyl)-8-quinolinyl ester-1H-indole-3-carboxylic
 14 acid).
 15 (FFFF) Benocyclidine (Other names include: BCP, BTCP, and
 16 Benzothiophenylcyclohexylpiperidine).
 17 (GGGG) 25B-NBOMe (Other names include: 2C-B-NBOMe
 18 and 4-Bromo-2,
 19 5-dimethoxy-N-[(2-Methoxyphenyl)methyl]
 20 benzeneethanamine).
 21 (HHHH) APB (Other names include: (2-Aminopropyl)
 22 Benzofuran).
 23 (IIII) AB-PINACA
 24 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
 25 indazole-3-carboxamide).
 26 (JJJJ) AB-FUBINACA
 27 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-
 28 1H-indazole-3-carboxamide).
 29 (KKKK) ADB-PINACA
 30 (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-in
 31 daole-3-carboxamide).
 32 (LLLL) Fluoro ADBICA (N-(1-Amino-3,3-
 33 dimethyl-1-oxobutan-2-yl)-(fluoropentyl)-1H-indole-3-
 34 carboxamide).
 35 (MMMM) APDB (Other names include: -EMA,
 36 -Desoxy-MDA, and (2-Aminopropyl)-2,3-
 37 dihydrobenzofuran).
 38 (NNNN) THJ-2201 (Other names include: AM2201 indazole
 39 analog, Fluoropentyl-JWH-018 indazole, and
 40 5-Fluoro-THJ-018).
 41 (OOOO) AM 2201 benzimidazole analog (Other names
 42 include: FUBIMINA, FTHJ, and (1-(5-fluoropentyl)-1H-



1 benzo[d]imidazol-2-yl)(naphthalene-1-yl)methanone).
 2 (PPPP) MN-25 (Other names include: 7-methoxy-1-
 3 [2-(4-morpholinyl)ethyl]-N-[1S, 2S, 4R]-1,3,3-
 4 trimethylbicyclo[2.2.1]hept-2-yl]-1H-indole-3-carboxamide
 5 and UR-12).
 6 (QQQQ) FUB-PB-22 (Other names include:
 7 Quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 8 (RRRR) FUD-PB-22 (Other names include:
 9 Naphthalen-1-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 10 (SSSS) 5-Fluoro-AB-PINACA (Other names include:
 11 AB-PINACA 5-fluoro analog and N-(1-amino-3-methyl-
 12 oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
 13 carboxamide).
 14 (TTTT) 4-MePPP (Other names include:
 15 4-methyl-alpha-pyrrolidinopropiophenone).
 16 (UUUU) alpha-PBP (Other names include:
 17 Alpha-pyrrolidinobutiophenone).
 18 (VVVV) AB-CHMINACA (Other names include:
 19 (N-[1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl-
 20 1)-1H-indazole-3-carboxamide).
 21 (WWWW) Acetyl fentanyl (Other names include:
 22 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide).
 23 (~~XXXX~~) (WWWW) M e x e d r o n e
 24 (3-methoxy-2-(methylamino)-1-(p-tolyl)propan-1-one).
 25 (XXXX) MT-45,
 26 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine).
 27 (YYYY) methyl 2-(1-(5-fluoropentyl)- 1H-indazole-3-
 28 carboxamido) -3,3-dimethylbutanoate [5F-ADB;
 29 5F-MDMB-PINACA].
 30 (ZZZZ) methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-
 31 carboxamido)-3-methylbutanoate [5F-AMB].
 32 (AAAAA) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)
 33 -1-(4-fluorobenzyl)- 1H-indazole-3-carboxamide
 34 [ADB-FUBINACA].
 35 (BBBBB) N-(adamantan-1-yl)-1-(5-fluoropentyl)-
 36 1H-indazole-3- carboxamide [5F-APINACA, 5F-AKB48].
 37 (CCCCC) methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
 38 carboxamido)-3,3-dimethylbutanoate [MDMB-CHMICA,
 39 MMB-CHMINACA].
 40 (DDDDD) methyl 2-(1-(4-fluorobenzyl)-
 41 1H-indazole-3-carboxamido)- 3,3-dimethylbutanoate
 42 [MDMB-FUBINACA].



- 1 (EEEE) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
 2 (cyclohexylmethyl)-1H-indazole-3-carboxamide
 3 [MAB-CHMINACA and ADB-CHMINACA].
 4 (FFFF) Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-
 5 3-carboxamido)-3-methylbutanoate [FUB-AMB,
 6 MMB-FUBINACA, AMB-FUBINACA].
 7 (GGGG) 3,4-dichloro-N-[(1dimethylamino)cyclohexylme
 8 thyl]benzamide [AH7921].
 9 (HHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1
 10 H-indole-3-carboxylate (trivial name: NM2201; CBL2201)
 11 (IIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1
 12 H-indazole-3-carboxamide (trivial name:
 13 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINA
 14 CA; 4-CN-CUMYL BINACA; CUMYL-4CN-BINACA;
 15 SGT-78).
 16 (JJJJ) methyl 2-(1-(cyclohexylmethyl)-1
 17 H-indole-3-carboxamido)-3-methylbutanoate (trivial
 18 names: MMB-CHMICA, AMB-CHMICA).
 19 (KKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1
 20 H-pyrrolo[2,3-b]pyridine-3-carboxamide (trivial name:
 21 5F-CUMYL-P7AICA).
 22 (LLLL) N-1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-p
 23 entanone (N-ethylpentylone, ephylone).
 24 (MMMM) Synthetic cathinone,
 25 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one
 26 (N-ethylpentylone, ephylone) and its optical, positional,
 27 and geometric isomers, salts, and salts of isomers.
 28 (2) Any compound structurally derived from
 29 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by
 30 substitution at the nitrogen atom of the indole ring by alkyl,
 31 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 32 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 33 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 34 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 35 or not further substituted in the indole ring to any extent and
 36 whether or not substituted in the naphthyl ring to any extent.
 37 (3) Any compound structurally derived from 3-(1-naphthoyl)
 38 pyrrole by substitution at the nitrogen atom of the pyrrole ring by
 39 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
 40 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,
 41 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl,
 42 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl



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



- 1 alkylendioxy, alkoxy, haloalkyl, hydroxyl, or halide
2 substituents, whether or not further substituted in the ring
3 system by one or more other univalent substituents;
4 (B) by substitution at the 3-position with an acyclic alkyl
5 substituent;
6 (C) by substitution at the 2-amino nitrogen atom with alkyl,
7 dialkyl, benzyl, or methoxybenzyl groups; or
8 (D) by inclusion of the 2-amino nitrogen atom in a cyclic
9 structure.
- 10 (9) Any compound structurally derived from 3-tetramethyl
11 cyclopropanoylindole with substitution at the nitrogen atom of the
12 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
13 cycloalkylmethyl, cycloalkylethyl,
14 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl) ethyl,
15 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
16 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
17 or not further substituted in the indole ring to any extent and
18 whether or not substituted in the tetramethylcyclopropyl ring to
19 any extent.
- 20 (10) Any compound containing a N-(1-adamantyl)-
21 1H-indazole-3-carboxamide structure with substitution at the
22 nitrogen atom of the indazole 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-methyl-2-pyrrolidinyl) methyl,
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 indazole 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 AKB48.
- 32 (11) Any compound containing a N-(1-adamantyl)-
33 1H-indole-3-carboxamide structure with substitution at the
34 nitrogen atom of the indole ring by an alkyl, haloalkyl,
35 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
36 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
37 1-(N-methyl-2-pyrrolidinyl) methyl,
38 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
39 group, whether or not further substituted at the nitrogen atom of
40 the carboxamide to any extent, whether or not further substituted
41 in the indole ring to any extent, and whether or not further
42 substituted on the adamantyl ring system to any extent. An



1 example of this structural class includes STS-135.

2 (12) Any compound containing a 3-(1-adamantoyl)indole
3 structure with substitution at the nitrogen atom of the indole ring
4 by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
5 cycloalkylethyl, 1-(N-methyl-2- piperidinyl)methyl, or
6 2-(4-morpholinyl)ethyl, 1-(N-methyl-2- pyrrolidinyl)methyl,
7 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
8 group, whether or not further substituted on the adamantyl ring
9 system to any extent. An example of this structural class includes
10 AM-1248.

11 (13) Any compound determined to be a synthetic drug by rule
12 adopted under IC 25-26-13-4.1.

13 SECTION 3. IC 35-48-1-16.6 IS ADDED TO THE INDIANA
14 CODE AS A NEW SECTION TO READ AS FOLLOWS
15 [EFFECTIVE JULY 1, 2019]: **Sec. 16.6. "Fentanyl related
16 substance" means any substance not listed in schedule I through V
17 of IC 35-48-2 that is structurally related to fentanyl by one (1) or
18 more of the following modifications:**

19 **(1) Replacement of the phenyl portion of the phenethyl group
20 by any monocycle, whether or not further substituted in or on
21 the monocycle.**

22 **(2) Substitution in or on the phenethyl group with alkyl,
23 alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino, or nitro
24 groups.**

25 **(3) Substitution in or on the piperidine ring with alkyl,
26 alkenyl, alkoxy, ester, ether, hydroxyl, halo, haloalkyl, amino,
27 or nitro groups.**

28 **(4) Replacement of the aniline ring with any aromatic
29 monocycle whether or not further substituted in or on the
30 aromatic monocycle.**

31 **(5) Replacement of the N-propionyl group by another acyl
32 group.**

33 SECTION 4. IC 35-48-2-4, AS AMENDED BY P.L.89-2018,
34 SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
35 JULY 1, 2019]: Sec. 4. (a) The controlled substances listed in this
36 section are included in schedule I.

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

42 **4-fluoroisobutyryl fentanyl**



- 1 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-
2 piperidinyl]-N-phenylacetamide) (9815)
- 3 **Acetyl fentanyl (Other names include:**
4 **N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide)**
- 5 Acetylmethadol (9601)
- 6 Acrylfentanyl. Other name: N-(1-phenethylpiperidin-4-yl)-
7 N-phenylacrylamide
- 8 Allylprodine (9602)
- 9 Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
10 thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)
- 11 Alphacetylmethadol (9603)
- 12 Alphameprodine (9604)
- 13 Alphamethadol (9605)
- 14 Alphamethylfentanyl (9814)
- 15 Benzethidine (9606)
- 16 Beta-hydroxy-3-methylfentanyl (9831). Other name:
17 N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl
18]-N-phenylpropanamide
- 19 Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-
20 phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)
- 21 Betacetylmethadol (9607)
- 22 Betameprodine (9608)
- 23 Betamethadol (9609)
- 24 Betaprodine (9611)
- 25 Clonitazene (9612)
- 26 Cyclopentyl fentanyl. Other name:
27 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide
- 28 Dextromoramide (9613)
- 29 Diampromide (9615)
- 30 Diethylthiambutene (9616)
- 31 Difenoxin (9168)
- 32 Dimenoxadol (9617)
- 33 Dimepheptanol (9618)
- 34 Dimethylthiambutene (9619)
- 35 Dioxaphetyl butyrate (9621)
- 36 Dipipanone (9622)
- 37 Ethylmethylthiambutene (9623)
- 38 Etonitazene (9624)
- 39 Etoxidine (9625)
- 40 **Fentanyl related substances.**
- 41 **Furanyl fentanyl.**
- 42 Furethidine (9626)



1	Hydroxypethidine (9627)
2	Isobutyryl fentanyl. Other name:
3	N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide
4	Ketobemidone (9628)
5	Levomoramide (9629)
6	Levophenacymorphan (9631)
7	Methoxyacetyl fentanyl. Other name:
8	2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide
9	3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-
10	piperidyl]-N-phenyl-propanamide](9813)
11	3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
12	piperidinyl]-N-phenylpropanamide) (9833)
13	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
14	Morpheridine (9632)
15	N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl),
16	including any isomers, salts, or salts of isomers (9818)
17	N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl] 25 piperidin-4-yl]-
18	N-phenylpropionamide, also known as N-[1-[2-hydroxy-2-
19	(2-thienyl)ethyl] -4- piperidinyl]- N-phenylpropanamide,
20	(beta-hydroxythiofentanyl)
21	N-(4-chlorophenyl)- N-(1-phenethylpiperidin-4-yl)
22	isobutyramide (para-chloroisobutyryl fentanyl)
23	N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)
24	acetamide (ocfentanil)
25	N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
26	(para-fluorobutyryl fentanyl)
27	N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also
28	known as N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide,
29	(butyryl fentanyl)
30	N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (valeryl
31	fentanyl)
32	N-(4-methoxyphenyl)-N-(1-phenethylpiperidin -4-yl)
33	butyramide (para-methoxybutyryl fentanyl)
34	N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide
35	(thenylfentanyl), including any isomers, salts, or salts of isomers
36	(9834)
37	N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide
38	(isobutyryl fentanyl)
39	N - (1 - p h e n e t h y l p i p e r i d i n - 4 - y l) -
40	Nphenylcyclopentanecarboxamide (cyclopentyl fentanyl)
41	Noracymethadol (9633)
42	Norlevorphanol (9634)



- 1 Normethadone (9635)
 2 Norpipanone (9636)
 3 O c f e n t a n i l . O t h e r n a m e :
 4 N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)
 5 acetamide
 6 Ortho-fluorofentanyl or 2-fluorofentanyl. Other name:
 7 N-(2-fluorophenyl)-N-
 8 (1-phenethylpiperidin-4-yl)propionamide
 9 Para-chloroisobutyryl fentanyl. Other name:
 10 N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide
 11 Para-fluorobutyryl fentanyl. Other name:
 12 N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
 13 Para-fluorofentanyl (N-(4-fluorophenyl)-N-
 14 [1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
 15 Para-methoxybutyryl fentanyl. Other name:
 16 N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
 17 Phenadoxone (9637)
 18 Phenampromide (9638)
 19 Phenomorphan (9647)
 20 Phenoperidine (9641)
 21 PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
 22 Piritramide (9642)
 23 Proheptazine (9643)
 24 Properidine (9644)
 25 Propiram (9649)
 26 Racemoramide (9645)
 27 Tetrahydrofuranyl fentanyl. Other name:
 28 N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carb
 29 oxamide
 30 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
 31 piperidinyl]-propanamide) (9835)
 32 Tilidine (9750)
 33 Trimeperidine (9646)
 34 U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-
 35 N-methyl- benzamide)
 36 V a l e r y l f e n t a n y l . O t h e r n a m e :
 37 N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide
 38 (c) Opium derivatives. Any of the following opium derivatives, their
 39 salts, isomers, and salts of isomers, unless specifically excepted by rule
 40 of the board or unless listed in another schedule, whenever the
 41 existence of these salts, isomers, and salts of isomers is possible within
 42 the specific chemical designation:



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



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



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



- 1 (B) the seeds of the plant;
 2 (C) any extract from any part of the plant; and
 3 (D) every compound, manufacture, salt, derivative, mixture, or
 4 preparation of the plant, its seeds, or extracts.
- 5 (36) 5-Methoxy-N,N-Dimethyltryptamine. Some trade or other
 6 names: 5-methoxy-3-[2- (dimethylamino)ethyl]indole;
 7 5-MeO-DMT.
 8 (37) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E).
 9 (38) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D).
 10 (39) 2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C).
 11 (40) 2-(4-Iodo-2,5-dimethoxyphenyl) ethanamine (2C-I).
 12 (41) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl] ethanamine (2C-T-2).
 13 (42) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl] ethanamine
 14 (2C-T-4).
 15 (43) 2-(2,5-Dimethoxyphenyl) ethanamine (2C-H).
 16 (44) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N).
 17 (45) 2-(2,5-Dimethoxy-4-(n)-propylphenyl) ethanamine (2C-P).
 18 (46) Deschloroketamine (2-Phenyl-2-
 19 (methylamino)cyclohexanone).
 20 (47) 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-
 21 ethyltryptamine).
 22 (48) N-methyltryptamine (1H-Indole-3-ethanamine, N-methyl-).
 23 (e) Depressants. Unless specifically excepted in a rule adopted by
 24 the board or unless listed in another schedule, any material, compound,
 25 mixture, or preparation which contains any quantity of the following
 26 substances having a depressant effect on the central nervous system,
 27 including its salts, isomers, and salts of isomers whenever the existence
 28 of such salts, isomers, and salts of isomers is possible within the
 29 specific chemical designation:
 30 Etizolam (4-(2- chlorophenyl)-2- ethyl-9- methyl- 6H-
 31 thieno[3,2-f] [1,2,4] triazolo[4,3-a] [1,4]diazepine) (other names
 32 include: Etilaam, Etizest, Depas, Etizola, Sedekopan, and
 33 Pasaden)
 34 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-
 35 4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine)
 36 Gamma-hydroxybutyric acid (other names include GHB;
 37 gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium
 38 oxybate; sodium oxybutyrate) (2010)
 39 Mecloqualone (2572)
 40 Methaqualone (2565)
 41 (f) Stimulants. Unless specifically excepted or unless listed in
 42 another schedule, any material, compound, mixture, or preparation that



1 contains any quantity of the following substances having a stimulant
 2 effect on the central nervous system, including its salts, isomers, and
 3 salts of isomers:

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

6 Aminorex (1585). Other names: aminoxaphen;
 7 2-amino-5-phenyl-2-oxazoline; or
 8 4,5-dihydro-5-phenyl-2-oxazolamine.

9 Cathinone (1235). Some trade or other names:
 10 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone;
 11 2-aminopropiophenone; and norephedrone.

12 Fenethylamine (1503).

13 N-Benzylpiperazine (7493). Other names: BZP; and
 14 1-benzylpiperazine.

15 N-ethylamphetamine (1475).

16 Methcathinone (1237). Some other trade names:
 17 2-Methylamino-1-Phenylpropan-I-one; Ephedrone;
 18 Monomethylpropion; UR 1431.

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

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

23 SECTION 5. IC 35-48-2-6, AS AMENDED BY P.L.89-2018,
 24 SECTION 3, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 25 JULY 1, 2019]: Sec. 6. (a) The controlled substances listed in this
 26 section are included in schedule II.

27 (b) Any of the following substances, except those narcotic drugs
 28 listed in other schedules, whether produced directly or indirectly by
 29 extraction from substances of vegetable origin, or independently by
 30 means of chemical synthesis, or by combination of extraction and
 31 chemical synthesis:

32 (1) Opium and opiate, and any salt, compound, derivative, or
 33 preparation of opium or opiate, excluding apomorphine,
 34 dextrorphan, nalbuphine, naloxone, naltrexone, and their
 35 respective salts but including:

36 (A) raw opium (9600);

37 (B) opium extracts (9610);

38 (C) opium fluid extracts (9620);

39 (D) powdered opium (9639);

40 (E) granulated opium (9640);

41 (F) tincture of opium (9630);

42 (G) codeine (9050);



- 1 (H) dihydroetorphine (9334);
 2 (I) ethylmorphine (9190);
 3 (J) etorphine hydrochloride (9059);
 4 (K) hydrocodone (9193), and any hydrocodone combination
 5 product, as determined by the federal Food and Drug
 6 Administration;
 7 (L) hydromorphone (9150);
 8 (M) metopon (9260);
 9 (N) morphine (9300);
 10 (O) oxycodone (9143);
 11 (P) oxymorphone (9652);
 12 (Q) thebaine (9333); and
 13 (R) oripavine.
- 14 (2) Any salt, compound, isomer, derivative, or preparation thereof
 15 which is chemically equivalent or identical with any of the
 16 substances referred to in subdivision (b)(1) of this section, but not
 17 including the isoquinoline alkaloids of opium.
- 18 (3) Opium poppy and poppy straw.
- 19 (4) Cocaine (9041).
- 20 (5) Concentrate of poppy straw (the crude extract of poppy straw
 21 in either liquid, solid, or powder form which contains the
 22 phenanthrene alkaloids of the opium poppy) (9670).
- 23 (c) Opiates. Any of the following opiates, including their isomers,
 24 esters, ethers, salts, and salts of isomers, esters, and ethers whenever
 25 the existence of these isomers, esters, ethers, and salts is possible
 26 within the specific chemical designation:
- 27 Alfentanil (9737)
 28 Alphaprodine (9010)
 29 Anileridine (9020)
 30 Bezitramide (9800)
 31 Bulk dextropropoxyphene (nondosage forms) (9273)
 32 Carfentanil (9743)
 33 Dihydrocodeine (9120)
 34 Diphenoxylate (9170)
 35 Fentanyl (9801)
 36 Isomethadone (9226)
 37 Levo-alpha-acetylmethadol (9648). Other names:
 38 Levo-alpha-acetylmethadol; levomethadyl acetate; and LAAM.
 39 Levomethorphan (9210)
 40 Levorphanol (9220)
 41 Metazocine (9240)
 42 Methadone (9250)



- 1 Methadone-Intermediate, 4-cyano-2-dimethyl-amino-4,
 2 4-diphenyl butane (9254)
 3 Moramide-Intermediate, 2-methyl-3-morpholino-1,
 4 1-diphenylpropane- carboxylic acid (9802)
 5 Pethidine (Meperidine) (9230)
 6 Pethidine-Intermediate- A, 4-cyano-1-methyl-4-phenylpiperidine
 7 (9232)
 8 Pethidine-Intermediate-B,
 9 ethyl-4-phenylpiperidine-4-carboxylate (9233)
 10 Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carbo
 11 xylic acid (9234)
 12 Phenazodine (9715)
 13 Piminodine (9730)
 14 Racemethorphan (9732)
 15 Racemorphan (9733)
 16 Remifentanil (9739)
 17 Sufentanil (9740)
 18 Tapentadol
 19 **Thiafentanil**
 20 (d) Stimulants. Any material compound, mixture, or preparation
 21 which contains any quantity of the following substances having a
 22 potential for abuse associated with a stimulant effect on the central
 23 nervous system:
 24 (1) Amphetamine, its salts, optical isomers, and salts of its optical
 25 isomers (1100).
 26 (2) Methamphetamine, including its salts, isomers, and salts of its
 27 isomers (1105).
 28 (3) Phenmetrazine and its salts (1631).
 29 (4) Methylphenidate (1724).
 30 (5) Lisdexamfetamine, its salts, its isomers, and salts of its
 31 isomers.
 32 (e) Depressants. Unless specifically excepted by rule of the board
 33 or unless listed in another schedule, any material, compound, mixture,
 34 or preparation which contains any quantity of the following substances
 35 having a depressant effect on the central nervous system, including its
 36 salts, isomers, and salts of isomers whenever the existence of such
 37 salts, isomers, and salts of isomers is possible within the specific
 38 chemical designation:
 39 Amobarbital (2125)
 40 Glutethimide (2550)
 41 Pentobarbital (2270)
 42 Phencyclidine (7471)



- 1 Secobarbital (2315)
- 2 (f) Immediate precursors. Unless specifically excepted by rule of the
- 3 board or unless listed in another schedule, any material, compound,
- 4 mixture, or preparation which contains any quantity of the following
- 5 substances:
- 6 (1) Immediate precursor to amphetamine and methamphetamine:
- 7 Phenylacetone (8501). Some trade or other names:
- 8 phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl
- 9 ketone.
- 10 (2) Immediate precursors to phencyclidine (PCP):
- 11 (A) 1-phenylcyclohexylamine (7460); or
- 12 (B) 1-piperidinocyclohexanecarbonitrile (PCC) (8603).
- 13 (3) Immediate precursor to fentanyl:
- 14 4-Anilino-N-Phenethyl-4-Piperidine (ANPP).
- 15 (g) Hallucinogenic substances:
- 16 Dronabinol **oral solution**. Other name:
- 17 (-)-delta-9-trans-tetrahydrocannabinol (delta-9-THC).
- 18 Nabilone (7379). Other name: (+/-)-trans-3-
- 19 (1,1-dimethylheptyl)-6, 6a, 7, 8, 10, 10a-hexahydro-1-hydroxy-6,
- 20 6-dimethyl-9H-dibenzo [b,d] pyran-9-one.
- 21 SECTION 6. IC 35-48-2-12, AS AMENDED BY P.L.283-2013,
- 22 SECTION 4, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
- 23 JULY 1, 2019]: Sec. 12. (a) The controlled substances listed in this
- 24 section are included in schedule V.
- 25 (b) Narcotic drugs containing nonnarcotic active medicinal
- 26 ingredients. Any compound, mixture, or preparation containing any of
- 27 the following narcotic drugs, or their salts calculated as the free
- 28 anhydrous base or alkaloid, in the following quantities, which shall
- 29 include one (1) or more nonnarcotic active medicinal ingredients in
- 30 sufficient proportion to confer upon the compound, mixture, or
- 31 preparation, valuable medicinal qualities other than those possessed by
- 32 the narcotic drug alone:
- 33 (1) Not more than 200 milligrams of codeine per 100 milliliters
- 34 or per 100 grams.
- 35 (2) Not more than 100 milligrams of dihydrocodeine per 100
- 36 milliliters or per 100 grams.
- 37 (3) Not more than 100 milligrams of ethylmorphine per 100
- 38 milliliters or per 100 grams.
- 39 (4) Not more than 2.5 milligrams of diphenoxylate and not less
- 40 than 25 micrograms of atropine sulfate per dosage unit.
- 41 (5) Not more than 100 milligrams of opium per 100 milliliters or
- 42 per 100 grams.



- 1 (6) Not more than 0.5 milligrams of difenoxin (9168), and not less
2 than 25 micrograms of atropine sulfate per dosage unit.
3 (c) **Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]**
4 **butanamide)**.
5 (d) Pregabalin (2782).
6 ~~(d)~~ (e) Pyrovalerone (1485).
7 ~~(e)~~ (f) Lacosamide [(R)-2-acetoamido-N-benzyl-
8 3-methoxy-propionamide].
9 (g) Epidiolex.



COMMITTEE REPORT

Madam President: The Senate Committee on Rules and Legislative Procedure, to which was referred Senate Bill No. 631, has had the same under consideration and begs leave to report the same back to the Senate with the recommendation that said bill be AMENDED as follows:

Delete everything after the enacting clause and insert the following:

(SEE TEXT OF BILL)

and when so amended that said bill be reassigned to the Senate Committee on Corrections and Criminal Law.

(Reference is to SB 631 as introduced.)

BRAY, Chairperson

COMMITTEE REPORT

Madam President: The Senate Committee on Corrections and Criminal Law, to which was referred Senate Bill No. 631, has had the same under consideration and begs leave to report the same back to the Senate with the recommendation that said bill be AMENDED as follows:

Page 1, between the enacting clause and line 1, begin a new paragraph and insert:

"SECTION 1. IC 35-31.5-2-130.5 IS ADDED TO THE INDIANA CODE AS A **NEW** SECTION TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2019]: **Sec. 130.5. "Fentanyl related substance", for purposes of IC 35-48, has the meaning set forth in IC 35-48-1-16.6.**"

Page 6, strike lines 16 through 17.

Page 6, line 18, strike "(XXXX)" and insert "(WWWW)".

Page 6, line 20, delete "(YYYY)" and insert "(XXXX)".

Page 6, delete lines 22 through 37.

Page 6, line 38, delete "(HHHHH)" and insert "(YYYY)".

Page 6, line 41, delete "(IIII)" and insert "(ZZZZ)".

Page 7, line 1, delete "(JJJJ)" and insert "(AAAAA)".

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Page 7, line 4, delete "(KKKKK)" and insert "(BBBBB)".

Page 7, line 6, delete "(LLLLL)" and insert "(CCCCC)".

Page 7, line 9, delete "(MMMMM)" and insert "(DDDDD)".

Page 7, line 12, delete "(NNNNN)" and insert "(EEEEEE)".

Page 7, line 15, delete "(OOOOO)" and insert "(FFFFFF)".

Page 7, line 18, delete "(PPPPP)" and insert "(GGGGG)".

Page 7, delete lines 20 through 28.

Page 7, line 29, delete "(TTTTT)" and insert "(HHHHH)".

Page 7, line 31, delete "(UUUUU)" and insert "(IIIII)".

Page 7, line 36, delete "(VVVVV)" and insert "(JJJJJ)".

Page 7, line 39, delete "(WWWWW)" and insert "(KKKKK)".

Page 8, line 1, delete "(XXXXX)" and insert "(LLLLL)".

Page 8, line 3, delete "(YYYYY)" and insert "(MMMMM)".

Page 10, between lines 33 and 34, begin a new paragraph and insert:
 "SECTION 4. IC 35-48-1-16.6 IS ADDED TO THE INDIANA CODE AS A NEW SECTION TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2019]: **Sec. 16.6. "Fentanyl related substance" means any substance not listed in schedule I through V of IC 35-48-2 that is structurally related to fentanyl by one (1) or more of the following modifications:**

(1) Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle.

(2) Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino, or nitro groups.

(3) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxy, ester, ether, hydroxyl, halo, haloalkyl, amino, or nitro groups.

(4) Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle.

(5) Replacement of the N-propionyl group by another acyl group.

SECTION 5. IC 35-48-2-4, AS AMENDED BY P.L.89-2018, SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2019]: 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:

4-fluoroisobutyryl fentanyl

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

Acetyl fentanyl (Other names include: N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide)

Acetylmethadol (9601)

Acrylfentanyl. Other name: N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide

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)

Cyclopentyl fentanyl. Other name: N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide

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)

Fentanyl related substances.



Furanyl fentanyl.

Furethidine (9626)

Hydroxypethidine (9627)

Isobutyryl fentanyl. Other name:

N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide

Ketobemidone (9628)

Levomoramide (9629)

Levophenacymorphan (9631)

Methoxyacetyl fentanyl. Other name:

2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide

3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenyl-propanamide](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-hydroxy-2-(thiophen-2-yl)ethyl] 25 piperidin-4-yl]-N-phenylpropionamide, also known as N-[1-[2-hydroxy-2-(2-thienyl)ethyl] -4- piperidinyl]- N-phenylpropanamide, (beta-hydroxythiofentanyl)**N-(4-chlorophenyl)- N-(1-phenethylpiperidin-4-yl) isobutyramide (para-chloroisobutyryl fentanyl)****N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl) acetamide (ocfentanil)****N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl) butyramide (para-fluorobutyryl fentanyl)****N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl)****N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (valeryl fentanyl)****N-(4-methoxyphenyl)-N-(1-phenethylpiperidin -4-yl) butyramide (para-methoxybutyryl fentanyl)**

N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), including any isomers, salts, or salts of isomers (9834)

N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl)**N - (1 - p h e n e t h y l p i p e r i d i n - 4 - y l) - Nphenylcyclopentanecarboxamide (cyclopentyl fentanyl)**

Noracymethadol (9633)
 Norlevorphanol (9634)
 Normethadone (9635)
 Norpipanone (9636)
 O c f e n t a n i l . O t h e r n a m e :
 N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)
 acetamide
 Ortho-fluorofentanyl or 2-fluorofentanyl. Other name:
 N-(2-fluorophenyl)-N-
 (1-phenethylpiperidin-4-yl)propionamide
 Para-chloroisobutyryl fentanyl. Other name:
 N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide
 Para-fluorobutyryl fentanyl. Other name:
 N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
 Para-fluorofentanyl (N-(4-fluorophenyl)-N-
 [1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
 Para-methoxybutyryl fentanyl. Other name:
 N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
 Phenadoxone (9637)
 Phenampromide (9638)
 Phenomorphan (9647)
 Phenoperidine (9641)
 PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
 Piritramide (9642)
 Proheptazine (9643)
 Properidine (9644)
 Propiram (9649)
 Racemoramide (9645)
 Tetrahydrofuranyl fentanyl. Other name:
 N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carb
 oxamide
 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)
 V a l e r y l f e n t a n y l . O t h e r n a m e :
 N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide

(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)
 Hydromorphanol (9301)
 Methyl-desorphine (9302)
 Methyl-dihydromorphine (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):

- (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- α -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).
- (46) Deschloroketamine (2-Phenyl-2-(methylamino)cyclohexanone).
- (47) 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-ethyltryptamine).
- (48) N-methyltryptamine (1H-Indole-3-ethanamine, N-methyl-).
- (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][1,4]diazepine) (other names include: Etilaam, Etizest, Depas, Etizola, Sedekopan, and Pasaden)
- Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine)
- 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."

Page 12, between lines 29 and 30, begin a new paragraph and insert: "**Thiafentanil**".

Page 14, line 13, delete "Pregabalin (2782)." and insert "**Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide)**".

Pregabalin (2782)."

Renumber all SECTIONS consecutively.

and when so amended that said bill do pass.

(Reference is to SB 631 as printed January 17, 2019.)

YOUNG M, Chairperson

Committee Vote: Yeas 8, Nays 1.



COMMITTEE REPORT

Mr. Speaker: Your Committee on Courts and Criminal Code, to which was referred Senate Bill 631, 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 20, line 20, beginning with "Thiafentaniil" begin a new line block indented.

Page 22, line 6, delete "Pregabalin" and insert "**(d)** Pregabalin".

Page 22, line 7, strike "(d)" and insert "**(e)**".

Page 22, line 8, strike "(e)" and insert "**(f)**".

Page 22, line 10, delete "(f)" and insert "**(g)**".

and when so amended that said bill do pass.

(Reference is to SB 631 as printed February 8, 2019.)

MCNAMARA

Committee Vote: yeas 11, nays 0.

