



Reprinted
January 28, 2022

SENATE BILL No. 209

DIGEST OF SB 209 (Updated January 27, 2022 3:44 pm - DI 106)

Citations Affected: IC 15-15; IC 24-4; IC 35-31.5; IC 35-48.

Synopsis: Drug schedules. Adds specified substances to the scheduled list of controlled substances. Replaces references to "delta-9 THC" with "THC". Provides a defense to certain controlled substance offenses if: (1) the controlled substance is hemp (as defined by federal law); and (2) the hemp was shipped in continuous transport from a licensed producer in another state to a licensed handler in another state.

Effective: July 1, 2022.

Young M, Brown L

January 6, 2022, read first time and referred to Committee on Corrections and Criminal Law.

January 20, 2022, amended, reported favorably — Do Pass.

January 24, 2022, read second time, amended, ordered engrossed.

January 25, 2022, engrossed. Returned to second reading.

January 27, 2022, re-read second time, amended, ordered engrossed.

SB 209—LS 6921/DI 92



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Second Regular Session of the 122nd General Assembly (2022)

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 2021 Regular Session of the General Assembly.

SENATE BILL No. 209

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 15-15-13-6.5, AS ADDED BY P.L.190-2019,
2 SECTION 7, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
3 JULY 1, 2022]: Sec. 6.5. As used in this chapter, "hemp product"
4 means a product derived from, or made by, processing hemp plants or
5 plant parts including derivatives, extracts, cannabinoids, isomers, acids,
6 salts, and salts of isomers. However, the term does not include:

7 (1) smokable hemp (as defined by IC 35-48-1-26.6); or
8 (2) products that contain a total ~~delta-9-tetrahydrocannabinol~~
9 **tetrahydrocannabinol** concentration of more than three-tenths
10 of one percent (0.3%) by weight.

11 SECTION 2. IC 24-4-21-3, AS AMENDED BY P.L.190-2019,
12 SECTION 24, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
13 JULY 1, 2022]: Sec. 3. A person may distribute low THC hemp extract
14 in Indiana only if the distributor has a certificate of analysis prepared
15 by an independent testing laboratory showing:

16 (1) that the low THC hemp extract is the product of a batch tested
17 by the independent testing laboratory;

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1 (2) that the independent testing laboratory determined that the
 2 batch contained not more than three-tenths percent (0.3%) total
 3 ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** (THC),
 4 including precursors **and isomers**, by weight, based on the testing
 5 of a random sample of the batch; and

6 (3) the cannabidiol percent present of the low THC hemp extract.

7 SECTION 3. IC 24-4-21-4, AS ADDED BY P.L.153-2018,
 8 SECTION 9, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 9 JULY 1, 2022]: Sec. 4. (a) Except as provided in subsection (b), low
 10 THC hemp extract must be distributed in packaging that contains the
 11 following information:

12 (1) A scannable bar code or QR code linked to a document that
 13 contains information with respect to the manufacture of the low
 14 THC hemp extract, including the:

15 (A) batch identification number;

16 (B) product name;

17 (C) batch date;

18 (D) expiration date, which must be not more than two (2) years
 19 from the date of manufacture;

20 (E) batch size;

21 (F) total quantity produced;

22 (G) ingredients used, including the:

23 (i) ingredient name;

24 (ii) name of the company that manufactured the ingredient;

25 (iii) company or product identification number or code, if
 26 applicable; and

27 (iv) ingredient lot number; and

28 (H) download link for a certificate of analysis for the low THC
 29 hemp extract.

30 (2) The batch number.

31 (3) The Internet address of a web site to obtain batch information.

32 (4) The expiration date.

33 (5) The number of milligrams of low THC hemp extract.

34 (6) The manufacturer.

35 (7) The fact that the product contains not more than three-tenths
 36 percent (0.3%) total ~~delta-9-tetrahydrocannabinol~~
 37 **tetrahydrocannabinol** (THC), including precursors **and isomers**,
 38 by weight.

39 (b) Before July 1, 2018, low THC hemp extract may be distributed
 40 in Indiana without having met the requirements described in subsection
 41 (a).

42 SECTION 4. IC 35-31.5-2-321, AS AMENDED BY P.L.61-2020,



1 SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
2 JULY 1, 2022]: Sec. 321. "Synthetic drug" means:

3 (1) a substance containing one (1) or more of the following
4 chemical compounds, including an analog of the compound **and**
5 **its isomers, salts, and salts of isomers:**

- 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 (naphthalen-1-yl-(1-butylandol-3-yl)methanone).
11 (E) JWH-081 (4-methoxynaphthalen-1-yl-(1-pentylindol-
12 3-yl)methanone).
13 (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
14 (G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-
15 naphthalen-1-yl-methanone).
16 (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
17 (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
18 (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
19 (K) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-
20 3-(2-methyloctan-2-yl)-
21 6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
22 (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-
23 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo
24 [c]chromen-1-ol).
25 (M) HU-308 ((1R,2R,5R)-2-[2,6-dimethoxy-4-
26 (2-methyloctan-2-yl)phenyl]-
27 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
28 (N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-
29 (1-methylethenyl)-2-cyclohexen-1-yl]-5
30 -pentyl-2,5-cyclohexadiene-1,4-dione).
31 (O) CP 55,940
32 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]-
33 5-(2-methyloctan-2-yl)phenol).
34 (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]-5-
35 (2-methyloctan-2-yl)phenol) and its homologues, or
36 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)
37 phenol, where side chain n=5, and homologues where side
38 chain n=4, 6, or 7.
39 (Q) WIN 55212-2
40 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
41 pyrrolo [1,2,3-de)-1,4-benzoxazin-
42 6-yl]-1-naphthalenylmethanone).



- 1 (R) RCS-4 ((4-methoxyphenyl)
 2 (1-pentyl-1H-indol-3-yl)methanone).
 3 (S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-
 4 indol-3-yl)-2-(2-methoxyphenyl)ethanone).
 5 (T) 4-Methylmethcathinone. Other name: mephedrone.
 6 (U) 3,4-Methylenedioxy-methcathinone. Other name:
 7 methylone.
 8 (V) Fluoromethcathinone.
 9 (W) 4-Methoxymethcathinone. Other name: methedrone.
 10 (X) 4-Ethylmethcathinone (4-EMC).
 11 (Y) Methylenedioxy-pyvalerone. Other name: MDPV.
 12 (Z) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
 13 (AA) JWH-098, or
 14 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
 15 (BB) JWH-164, or
 16 1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
 17 (CC) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
 18 (DD) JWH-201, or
 19 1-pentyl-3-(4-methoxyphenylacetyl)indole.
 20 (EE) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole.
 21 (FF) AM-694, or
 22 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
 23 (GG) CP 50,556-1, or
 24 [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylp
 25 ntan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1
 26 -yl] acetate.
 27 (HH) Dimethylheptylpyran, or DMHP.
 28 (II) 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP.
 29 (JJ) 6-APB [6-(2-aminopropyl)benzofuran].
 30 (LL) 7-hydroxymitragynine.
 31 (MM) α -PPP [α -pyrrolidinopropiophenone].
 32 (NN) α -PVP (desmethylpyrovalerone).
 33 (OO) AM-251.
 34 (PP) AM-1241.
 35 (QQ) AM-2201.
 36 (RR) AM-2233.
 37 (SS) Buphedrone (α -methylamino-butyrophenone (MABP)).
 38 (TT) Butylone.
 39 (UU) CP-47,497-C7.
 40 (VV) CP-47,497-C8.
 41 (WW) Desoxypipradol.
 42 (XX) Ethylone.



1	(YY) Eutylone.
2	(ZZ) Flephedrone.
3	(AAA) JWH-011.
4	(BBB) JWH-020.
5	(CCC) JWH-022.
6	(DDD) JWH-030.
7	(EEE) JWH-182.
8	(FFF) JWH-302.
9	(GGG) MDAI [5,6-methylenedioxy-2-aminoindane].
10	(HHH) Mitragynine.
11	(III) Naphyrone.
12	(JJJ) Pentedrone.
13	(LLL) Pentylone.
14	(MMM) Methoxetamine
15	[2-(3-methoxyphenyl)-2-(ethylamino)- cyclohexanone].
16	(NNN) A796,260 [1-(2-morpholin-4-ylethyl)-1H-indol-3-yl]-
17	(2,2,3,3-tetramethylcyclopropyl)methanone].
18	(OOO) AB-001[(1s,3s)-adamantan-1-yl)
19	(1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
20	(1-adamantoyl)indole].
21	(PPP) AM-356 [Methanandamide].
22	(QQQ) AM 1248 [1-[(1-methyl-2- piperidinyl) methyl]-
23	1H-indol-3-yl] tricyclo[3.3.1.1 ^{3,7}] dec-1-yl-methanone]or
24	[(1-[(N-methylpiperindin-2-yl)
25	Methyl]-3-(Adamant-1-oyl)indole].
26	(RRR) AM 2233 Azepane isomer [(2-iodophenyl)
27	(1-(1-methylazepan-3-yl)- 1H-indol-3-yl)methanone].
28	(SSS) CB-13 [1-Naphthalenyl
29	[4-(pentyoxy)- 1-naphthalenyl]methanone].
30	(TTT) UR-144 [(1-pentyl-1H-indol-3-yl)
31	(2,2,3,3-tetramethylcyclopropyl)-methanone].
32	(UUU) URB 597 [(3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl)-
33	cyclohexylcarbamate].
34	(VVV) URB602 [[1,1'-biphenyl]- 3-yl-carbamic acid,
35	cyclohexyl ester].
36	(WWW) URB 754 [6-methyl-2-[(4-methylphenyl)
37	amino]-1-benzoxazin-4-one].
38	(XXX) XLR-11 or 5-fluoro UR-144
39	(1-(5-fluoropentyl)-1H-indol-3-yl)
40	(2,2,3,3-tetramethylcyclopropyl)methanone].
41	(YYY) AKB48 (Other names include:
42	N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide;



1 1-pentyl-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-indazole-3-
2 carboxamide).
3 (ZZZ) 25I-NBOMe (Other names include:
4 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-
5 benzeneethanamine);
6 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
7 methyl]ethanamine).
8 (AAAA) 2C-C-NBOMe (Other names include: 25C-NBOMe;
9 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
10 methyl]ethanamine;
11 2,5-Dimethoxy-4-chloro-N-(2-methoxybenzyl)
12 phenethylamine).
13 (BBBB) 2NE-1 (Other names include: 1-Pentyl-3-
14 (1-adamantylamido)indole).
15 (CCCC) STS-135 (Other names include:
16 N-Adamantyl-1-fluoropentylindole-3- carboxamide
17 (1-5-fluoropentyl)-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-
18 indole-3-carboxamide).
19 (DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinyl
20 ester-1H-indole-2-carboxylic acid).
21 (EEEE) 5-Fluoro-PB-22 (Other names include:
22 1-(5-Fluoropentyl)-8-quinolinyl ester-1H-indole-3-carboxylic
23 acid).
24 (FFFF) Benocyclidine (Other names include: BCP, BTCP, and
25 Benzothiophenylcyclohexylpiperidine).
26 (GGGG) 25B-NBOMe (Other names include: 2C-B-NBOMe
27 and 4-Bromo-2,
28 5-dimethoxy-N-[(2-Methoxyphenyl)methyl]
29 benzeneethanamine).
30 (HHHH) APB (Other names include: (2-Aminopropyl)
31 Benzofuran).
32 (III) AB-PINACA
33 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
34 indazole-3-carboxamide).
35 (JJJ) AB-FUBINACA
36 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-
37 1H-indazole-3-carboxamide).
38 (KKKK) ADB-PINACA
39 (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
40 indazole-3-carboxamide).
41 (LLL) Fluoro ADBICA (N-(1-Amino-3,3-
42 dimethyl-1-oxobutan-2-yl)-(fluoropentyl)-1H-indole-3-



1 carboxamide).
 2 (MMMM) APDB (Other names include: -EMA,
 3 -Desoxy-MDA, and (2-Aminopropyl)-2,3-
 4 dihydrobenzofuran).
 5 (NNNN) THJ-2201 (Other names include: AM2201 indazole
 6 analog, Fluoropentyl-JWH-018 indazole, and
 7 5-Fluoro-THJ-018).
 8 (OOOO) AM 2201 benzimidazole analog (Other names
 9 include: FUBIMINA, FTHJ, and (1-(5-fluoropentyl)-1H-
 10 benzo[d]imidazol-2-yl)(naphthalene-1-yl)methanone).
 11 (PPPP) MN-25 (Other names include: 7-methoxy-1-
 12 [2-(4-morpholinyl)ethyl]-N-[1S, 2S, 4R]-1,3,3-
 13 trimethylbicyclo[2.2.1]hept-2-yl]-1H-indole-3-carboxamide
 14 and UR-12).
 15 (QQQQ) FUB-PB-22 (Other names include:
 16 Quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 17 (RRRR) FUD-PB-22 (Other names include:
 18 Naphthalen-1-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 19 (SSSS) 5-Fluoro-AB-PINACA (Other names include:
 20 AB-PINACA 5-fluoro analog and N-(1-amino-3-methyl-
 21 oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
 22 carboxamide).
 23 (TTTT) 4-MePPP (Other names include:
 24 4-methyl-alpha-pyrrolidinopropiophenone).
 25 (UUUU) alpha-PBP (Other names include:
 26 Alpha-pyrrolidinobutiophenone).
 27 (VVVV) AB-CHMINACA (Other names include:
 28 (N-[1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl-
 29 1)-1H-indazole-3-carboxamide).
 30 (WWWW) Mexedrone
 31 (3-methoxy-2-(methylamino)-1-(p-tolyl)propan-1-one).
 32 (XXXX) MT-45,
 33 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine).
 34 (YYYY) methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
 35 carboxamido)-3,3-dimethylbutanoate [5F-ADB;
 36 5F-MDMB-PINACA].
 37 (ZZZZ) methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
 38 carboxamido)-3-methylbutanoate [5F-AMB].
 39 (AAAAA) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-
 40 -1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
 41 [ADB-FUBINACA].
 42 (BBBBB) N-(adamantan-1-yl)-1-(5-fluoropentyl)-



- 1 1H-indazole-3- carboxamide [5F-APINACA, 5F-AKB48].
 2 (CCCCC) methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
 3 carboxamido)-3,3-dimethylbutanoate [MDMB-CHMICA,
 4 MMB-CHMINACA].
 5 (DDDDD) methyl 2-(1-(4-fluorobenzyl)-
 6 1H-indazole-3-carboxamido)- 3,3-dimethylbutanoate
 7 [MDMB-FUBINACA].
 8 (EEEEEE) N-(1-amino-3,3-dimethyl-1 -oxobutan-2-yl)-1-
 9 (cyclohexylmethyl)- 1H-indazole-3-carboxamide
 10 [MAB-CHMINACA and ADB-CHMINACA].
 11 (FFFFFF) Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-
 12 3-carboxamido)- 3-methylbutanoate [FUB-AMB,
 13 MMB-FUBINACA, AMB-FUBINACA].
 14 (GGGGG) 3,4-dichloro-N-[(1dimethylamino)cyclohexylme
 15 thyl]benzamide [AH7921].
 16 (HHHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1
 17 H-indole-3-carboxylate (trivial name: NM2201; CBL2201)
 18 (IIIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1
 19 H-indazole-3-carboxamide (trivial name:
 20 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINA CA;
 21 4-CN-CUMYL BINACA; CUMYL-4CN-BINACA; SGT-78).
 22 (JJJJJ) methyl 2-(1-(cyclohexylmethyl)-1
 23 H-indole-3-carboxamido)-3-methylbutanoate (trivial names:
 24 MMB-CHMICA, AMB-CHMICA).
 25 (KKKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1
 26 H-pyrrolo[2,3-b]pyridine-3-carboxamide (trivial name:
 27 5F-CUMYL-P7AICA).
 28 (LLLLL) N-1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-p
 29 entanone (N-ethylpentylone, ephylone).
 30 (MMMMM) Synthetic cathinone, 1-(1,3-benzodioxol-5-yl)-2-
 31 (ethylamino)- pentan-1-one (N-ethylpentylone, ephylone) and
 32 its optical, positional, and geometric isomers, salts, and salts
 33 of isomers.
 34 (NNNNN) ethyl
 35 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-
 36 dimethylbutanoate (trivial name: 5F-EDMB-PINACA).
 37 (OOOOO) methyl 2-(1-(5- fluoropentyl)-1H-indole-3-
 38 carboxamido)-3,3-dimethylbutanoate (trivial name:
 39 5F-MDMB-PICA).
 40 (PPPPP) N- (adamantan- 1-yl)- 1-(4-fluorobenzyl)-
 41 1H-indazole-3-carboxamide (trivial names: FUB-AKB48;
 42 FUB-APINACA; AKB48 N- (4-FLUOROBENZYL)).



- 1 (QQQQQ) 1-(5- fluoropentyl)-N-(2-phenylpropan-2-yl)-
 2 1H-indazole-3-carboxamide (trivial names:
 3 5F-CUMYL-PINACA; SGT-25).
 4 (RRRRR) (1-(4-fluorobenzyl)-1H-indol-3-
 5 yl)(2,2,3,3-tetramethylcyclopropyl) methanone (trivial name:
 6 FUB-144).
 7 (SSSSS) 4F-MDMB-BINACA.
 8 (TTTTT) N - e t h y l h e x e d r o n e
 9 (2-(ethylamino)-1-phenylhexan-1-one).
 10 (UUUUU) alpha-pyrrolidinohexanophenone
 11 (1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one.
 12 (VVVVV) alpha-pyrrolidinohexiophenone; trivial name:
 13 a-PHP.
 14 (WWWWW) 4'-methyl-alpha-pyrrolidinohexiophenone
 15 (1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one.
 16 (XXXXX) 4-methyl-alphaethylaminopentiophenone
 17 (2-(ethylamino)-1-(4-methylphenyl)pentan-1-one; trivial
 18 name: 4-MEAP.
 19 (YYYYY) 4'-methyl-alphapyrrolidinohexanophenone; trivial
 20 name: MPHP.
 21 (ZZZZZ) a l p h a p y r r o l i d i n o h e p t a p h e n o n e
 22 (1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one; trivial name: PV8.
 23 (AAAAA) 4'-chloro-alphapyrrolidinovalerophenone (1-(4-
 24 chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one.
 25 (BBBBB) 4'-chloro-alphapyrrolidinopentiophenone; trivial
 26 name: 4-chloro-a-PVP.
 27 (CCCCCC) **4,4'-dimethylaminorex (common name:**
 28 **4,4'-DMAR).**
 29 (2) Any compound structurally derived from
 30 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by
 31 substitution at the nitrogen atom of the indole ring by alkyl,
 32 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 33 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 34 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 35 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 36 or not further substituted in the indole ring to any extent and
 37 whether or not substituted in the naphthyl ring to any extent.
 38 (3) Any compound structurally derived from 3-(1-naphthoyl)
 39 pyrrole by substitution at the nitrogen atom of the pyrrole ring by
 40 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
 41 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,
 42 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl,



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



- 1 (A) by substitution in the ring system to any extent with alkyl,
 2 alkylenedioxy, alkoxy, haloalkyl, hydroxyl, or halide
 3 substituents, whether or not further substituted in the ring
 4 system by one or more other univalent substituents;
 5 (B) by substitution at the 3-position with an acyclic alkyl
 6 substituent;
 7 (C) by substitution at the 2-amino nitrogen atom with alkyl,
 8 dialkyl, benzyl, or methoxybenzyl groups; or
 9 (D) by inclusion of the 2-amino nitrogen atom in a cyclic
 10 structure.
- 11 (9) Any compound structurally derived from 3-tetramethyl
 12 cyclopropanoylindole with substitution at the nitrogen atom of the
 13 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
 14 cycloalkylmethyl, cycloalkylethyl,
 15 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl) ethyl,
 16 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
 17 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 18 or not further substituted in the indole ring to any extent and
 19 whether or not substituted in the tetramethylcyclopropyl ring to
 20 any extent.
- 21 (10) Any compound containing a N-(1-adamantyl)-
 22 1H-indazole-3-carboxamide structure with substitution at the
 23 nitrogen atom of the indazole ring by an alkyl, haloalkyl,
 24 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 25 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
 26 1-(N-methyl-2-pyrrolidinyl) methyl,
 27 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
 28 group, whether or not further substituted at the nitrogen atom of
 29 the carboxamide to any extent, whether or not further substituted
 30 in the indazole ring to any extent, and whether or not further
 31 substituted on the adamantyl ring system to any extent. An
 32 example of this structural class includes AKB48.
- 33 (11) Any compound containing a N-(1-adamantyl)-
 34 1H-indole-3-carboxamide structure with substitution at the
 35 nitrogen atom of the indole ring by an alkyl, haloalkyl,
 36 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 37 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
 38 1-(N-methyl-2-pyrrolidinyl) methyl,
 39 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
 40 group, whether or not further substituted at the nitrogen atom of
 41 the carboxamide to any extent, whether or not further substituted
 42 in the indole ring to any extent, and whether or not further



1 substituted on the adamantyl ring system to any extent. An
 2 example of this structural class includes STS-135.

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

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

14 SECTION 5. IC 35-48-1-17.4, AS ADDED BY P.L.61-2020,
 15 SECTION 3, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 16 JULY 1, 2022]: Sec. 17.4. (a) Except as provided in subsections (b)
 17 and (c), "isomer" means an optical isomer.

18 (b) "Isomer", as used in IC 35-48-2-4(d), means an optical,
 19 positional, or geometric isomer.

20 (c) "Isomer", as used in section 7 of this chapter, means an optical
 21 or geometric isomer. **"Isomer" means an optical, positional, or
 22 geometric isomer.**

23 SECTION 6. IC 35-48-1-17.5, AS AMENDED BY P.L.190-2019,
 24 SECTION 28, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 25 JULY 1, 2022]: Sec. 17.5. (a) "Low THC hemp extract" means a
 26 substance or compound that:

- 27 (1) is derived from or contains any part of the plant *Cannabis*
 28 *sativa* L. that meets the definition of hemp under IC 15-15-13-6;
- 29 (2) contains not more than three-tenths percent (0.3%) total
 30 ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** (THC),
 31 including precursors **and isomers**, by weight; and
- 32 (3) contains no other controlled substances.

33 (b) The term does not include:

- 34 (1) the harvested reproductive organ, whether immature or
 35 mature, of the female hemp plant; or
- 36 (2) smokable hemp.

37 SECTION 7. IC 35-48-2-4, AS AMENDED BY P.L.10-2021,
 38 SECTION 1, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
 39 JULY 1, 2022]: Sec. 4. (a) The controlled substances listed in this
 40 section are included in schedule I.

41 (b) Opiates. Any of the following opiates, including their isomers,
 42 esters, ethers, salts, and salts of isomers, esters, and ethers, unless



1 specifically excepted by rule of the board or unless listed in another
 2 schedule, whenever the existence of these isomers, esters, ethers, and
 3 salts is possible within the specific chemical designation:
 4 4-fluoroisobutyryl fentanyl
 5 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-
 6 piperidinyl]-N-phenylacetamide) (9815)
 7 Acetyl fentanyl (Other names include:
 8 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide)
 9 Acetylmethadol (9601)
 10 Acrylfentanyl. Other name: N-(1-phenethylpiperidin-4-yl)-
 11 N-phenylacrylamide
 12 Allylprodine (9602)
 13 Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
 14 thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)
 15 Alphacetylmethadol (9603)
 16 Alphameprodine (9604)
 17 Alphamethadol (9605)
 18 Alphamethylfentanyl (9814)
 19 Benzethidine (9606)
 20 Beta-hydroxy-3-methylfentanyl (9831). Other name:
 21 N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl
 22]-N-phenylpropanamide
 23 Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-
 24 phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)
 25 Betacetylmethadol (9607)
 26 Betameprodine (9608)
 27 Betamethadol (9609)
 28 Betaprodine (9611)
 29 **Brorphine**
 30 **2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1yl)-N,N-di**
 31 **ethylethan-1-amine (butonitazene)**
 32 Clonitazene (9612)
 33 Cyclopentyl fentanyl. Other name:
 34 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide
 35 Dextromoramide (9613)
 36 Diampromide (9615)
 37 Diethylthiambutene (9616)
 38 **N,N-diethyl-2-(2-(4-flourobenzyl)-5-nitro-1H-benzimidazol-**
 39 **1-yl)ethan-1-amine (flunitazene)**
 40 **N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)e**
 41 **than-1-amine (metodesnitazene)**
 42 **N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidaz**



1	ol-1-yl)ethan-1-amine (metonitazene)
2	N,N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine (protonitazene)
3	
4	Difenoxin (9168)
5	Dimenoxadol (9617)
6	Dimepheptanol (9618)
7	Dimethylthiambutene (9619)
8	Dioxaphetyl butyrate (9621)
9	Dipipanone (9622)
10	2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine (etodesnitazene; etazene)
11	
12	2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-benzimidazole (N-pyrrolidino etonitazene; etonitazepyne)
13	
14	Ethylmethylthiambutene (9623)
15	Etonitazene (9624)
16	Etoxidine (9625)
17	Fentanyl related substances.
18	Furanyl fentanyl.
19	Furethidine (9626)
20	Hydroxypethidine (9627)
21	Isobutyryl fentanyl. Other name:
22	N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide
23	Isotonitazene. Other name: N,N-diethyl-2-
24	(2-(4 isopropoxybenzyl)-5-nitro-1H-benzimidazol-
25	1-yl)ethan-1-amine)
26	Ketobemidone (9628)
27	Levomoramide (9629)
28	Levophenacymorphan (9631)
29	Methoxyacetyl fentanyl. Other name:
30	2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide
31	3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-
32	piperidyl]-N-phenyl-propanamide](9813)
33	3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
34	piperidinyl]-N-phenylpropanamide) (9833)
35	MDMB-4en-PINACA
36	4F-MDMB-BICA; 4-fluoro MDMB-BICA; 4F-MDMB-BUTICA;
37	Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,
38	3-dimethyl-butanoate
39	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
40	Morpheridine (9632)
41	N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl),
42	including any isomers, salts, or salts of isomers (9818)



1 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl] 25 piperidin-4-yl]-
2 N-phenylpropionamide, also known as N-[1-[2-hydroxy-2-
3 (2-thienyl)ethyl] -4- piperidinyl]- N-phenylpropanamide,
4 (beta-hydroxythiofentanyl)
5 N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl) isobutyramide
6 (para-chloroisobutyryl fentanyl)
7 N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)
8 acetamide (ocfentanil)
9 N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4 -yl) butyramide
10 (para-fluorobutyryl fentanyl)
11 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known
12 as N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl
13 fentanyl)
14 N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (valeryl
15 fentanyl)
16 N-(4-methoxyphenyl)-N-(1-phenethylpiperidin -4-yl) butyramide
17 (para-methoxybutyryl fentanyl)
18 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide
19 (thenylfentanyl), including any isomers, salts, or salts of isomers
20 (9834)
21 N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl
22 fentanyl)
23 N-(1-phenethylpiperidin-4-yl)-Nphenylcyclopentanecarboxamide
24 (cyclopentyl fentanyl)
25 Noracymethadol (9633)
26 Norlevorphanol (9634)
27 Normethadone (9635)
28 Norpipanone (9636)
29 O c f e n t a n i l . O t h e r n a m e :
30 N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)
31 acetamide
32 Ortho-fluorofentanyl or 2-fluorofentanyl. Other name:
33 N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide
34 Para-chloroisobutyryl fentanyl. Other name:
35 N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide
36 Para-fluorobutyryl fentanyl. Other name:
37 N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
38 Para-fluorofentanyl (N-(4-fluorophenyl)-N-
39 [1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
40 Para-methoxybutyryl fentanyl. Other name:
41 N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide
42 Phenadoxone (9637)



- 1 Phenampromide (9638)
 2 Phenomorphan (9647)
 3 Phenoperidine (9641)
 4 PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
 5 Pir tramide (9642)
 6 Proheptazine (9643)
 7 Properidine (9644)
 8 Propiram (9649)
 9 Racemoramide (9645)
 10 Tetrahydrofuranyl fentanyl. Other name:
 11 N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carb
 12 oxamide
 13 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
 14 piperidinyl]-propanamide) (9835)
 15 Tilidine (9750)
 16 Trimeperidine (9646)
 17 U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-
 18 N-methyl- benzamide)
 19 Valeryl fentanyl. Other name:
 20 N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide
 21 (c) Opium derivatives. Any of the following opium derivatives, their
 22 salts, isomers, and salts of isomers, unless specifically excepted by rule
 23 of the board or unless listed in another schedule, whenever the
 24 existence of these salts, isomers, and salts of isomers is possible within
 25 the specific chemical designation:
 26 Acetorphine (9319)
 27 Acetyldihydrocodeine (9051)
 28 Benzylmorphine (9052)
 29 Codeine methylbromide (9070)
 30 Codeine-N-Oxide (9053)
 31 Cyprenorphine (9054)
 32 Desomorphine (9055)
 33 Dihydromorphine (9145)
 34 Drotebanol (9335)
 35 Etorphine (except hydrochloride salt) (9056)
 36 Heroin (9200)
 37 Hydromorphenol (9301)
 38 Methyl desorphine (9302)
 39 Methyl dihydromorphine (9304)
 40 Morphine methylbromide (9305)
 41 Morphine methylsulfonate (9306)
 42 Morphine-N-Oxide (9307)



- 1 Myrophine (9308)
 2 Nicocodeine (9309)
 3 Nicomorphine (9312)
 4 Normorphine (9313)
 5 Pholcodine (9314)
 6 Thebacon (9315)
 7 (d) Hallucinogenic substances. Unless specifically excepted or
 8 unless listed in another schedule, any material, compound, mixture, or
 9 preparation which contains any quantity of the following
 10 hallucinogenic, psychedelic, or psychogenic substances, their salts,
 11 isomers, and salts of isomers whenever the existence of these salts,
 12 isomers, and salts of isomers is possible within the specific chemical
 13 designation (for purposes of this subsection only, the term "isomer"
 14 includes the optical, position, and geometric isomers): designation:
 15 (1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name:
 16 TCPy.
 17 (2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or
 18 other names: 4-Bromo-2, 5-Dimethoxy-a-methylphenethylamine;
 19 4-Bromo-2, 5-DMA.
 20 (3) 4-Bromo-2, 5-dimethoxyphenethylamine (7392). Some trade
 21 or other names:
 22 2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;
 23 alpha-desmethyl DOB; 2C-B, Nexus.
 24 (4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name:
 25 DOET.
 26 (5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348).
 27 Other name: 2C-T-7.
 28 (6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other
 29 names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.
 30 (7) 4-Methoxyamphetamine (7411). Some trade or other names:
 31 4-Methoxy-a-methylphenethylamine; Paramethoxyamphetamine;
 32 PMA.
 33 (8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401). Other
 34 Name: MMDA.
 35 (9) 5-Methoxy-N, N-diisopropyltryptamine, including any
 36 isomers, salts, or salts of isomers (7439). Other name:
 37 5-MeO-DIPT.
 38 (10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade
 39 and other names: 4-methyl-2,
 40 5-dimethoxy-a-methylphenethylamine; DOM; and STP.
 41 (11) 3, 4-methylenedioxy amphetamine (7400). Other name:
 42 MDA.



- 1 (12) 3,4-methylenedioxy-N-ethylamphetamine (7404). Other
 2 names: N-ethyl-alpha-methyl-3,4(methylenedioxy)
 3 phenethylamine; N-ethyl MDA; MDE; and MDEA.
 4 (13) 3, 4-methylenedioxymethamphetamine (MDMA) (7405).
 5 (14) 3, 4, 5-trimethoxy amphetamine (7390). Other name: TMA.
 6 (15) Alpha-ethyltryptamine (7249). Some trade and other names:
 7 Etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine;
 8 3-(2-aminobutyl) indole; [alpha]-ET; and AET.
 9 (16) Alpha-methyltryptamine (7432). Other name: AMT.
 10 (17) Bufotenine (7433). Some trade and other names:
 11 3-(B-Dimethylaminoethyl)-5-hydroxyindole;
 12 3-(2-dimethylaminonethyl)-5-indolol; N, N-dimethylserotonin;
 13 5-hydroxy-N, N-dimethyltryptamine; mappine.
 14 (18) Diethyltryptamine (7434). Some trade or other names: N,
 15 N-Diethyltryptamine; DET.
 16 (19) Dimethyltryptamine (7435). Some trade or other names:
 17 DMT.
 18 (20) Ibogaine (7260). Some trade and other names: 7-Ethyl-6, 6b,
 19 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6, 9-methano-5H-pyrido
 20 (1', 2': 1, 2, azepino 4, 5-b) indole; tabernanthe iboga.
 21 (21) Lysergic acid diethylamide (7315). Other name: LSD.
 22 (22) Marijuana (7360).
 23 (23) Mescaline (7381).
 24 (24) Parahexyl (7374). Some trade or other names:
 25 3-Hexyl-1-hydroxy-7, 8, 9, 10-Tetrahydro-6, 6,
 26 9-trimethyl-6H-dibenzo (b,d) pyran; Snyhexyl.
 27 (25) Peyote (7415), including:
 28 (A) all parts of the plant that are classified botanically as
 29 lophophora williamsii lemaire, whether growing or not;
 30 (B) the seeds thereof;
 31 (C) any extract from any part of the plant; and
 32 (D) every compound, manufacture, salt, derivative, mixture, or
 33 preparation of the plant, its seeds, or extracts.
 34 (26) N-ethyl-3-piperidyl benzilate (7482). Other name: DMZ.
 35 (27) N-hydroxy-3,4-methylenedioxyamphetamine (7402). Other
 36 names: N-hydroxy-alpha-methyl-3,4
 37 (methylenedioxy)phenethylamine; and N-hydroxy MDA.
 38 (28) N-methyl-3-piperidyl benzilate (7484). Other name: LBJ.
 39 (29) Psilocybin (7437).
 40 (30) Psilocyn (7438).
 41 (31) Tetrahydrocannabinols (7370), including synthetic
 42 equivalents of the substances contained in the plant, or in the



1 resinous extractives of Cannabis, sp. and synthetic substances,
 2 derivatives, and their isomers with similar chemical structure and
 3 pharmacological activity such as:

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

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

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

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

14 (32) Ethylamine analog of phencyclidine (7455). Some trade or
 15 other names: N-Ethyl-1-phenylcyclohexylamine;
 16 (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl)
 17 ethylamine; cyclohexamine; PCE.

18 (33) Pyrrolidine analog of phencyclidine (7458). Some trade or
 19 other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCP,_y; PHP.

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

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

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

26 (B) the seeds of the plant;

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

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

30 (36) 5-Methoxy-N,N-Dimethyltryptamine. Some trade or other
 31 names: 5-methoxy-3-[2- (dimethylamino)ethyl]indole;
 32 5-MeO-DMT.

33 (37) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E).

34 (38) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D).

35 (39) 2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C).

36 (40) 2-(4-Iodo-2,5-dimethoxyphenyl) ethanamine (2C-I).

37 (41) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl] ethanamine (2C-T-2).

38 (42) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl] ethanamine
 39 (2C-T-4).

40 (43) 2-(2,5-Dimethoxyphenyl) ethanamine (2C-H).

41 (44) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N).

42 (45) 2-(2,5-Dimethoxy-4-(n)-propylphenyl) ethanamine (2C-P).



- 1 (46) Deschloroketamine (2-Phenyl-2-
 2 (methylamino)cyclohexanone).
 3 (47) 4-Hydroxy-MET (4-Hydroxy-N-methyl-N-
 4 ethyltryptamine).
 5 (48) N-methyltryptamine (1H-Indole-3-ethanamine, N-methyl-).
 6 (e) Depressants. Unless specifically excepted in a rule adopted by
 7 the board or unless listed in another schedule, any material, compound,
 8 mixture, or preparation which contains any quantity of the following
 9 substances having a depressant effect on the central nervous system,
 10 including its salts, isomers, and salts of isomers whenever the existence
 11 of such salts, isomers, and salts of isomers is possible within the
 12 specific chemical designation:
 13 Etizolam (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-
 14 thieno[3,2-f][1,2,4]triazolo[4,3-a][1,4]diazepine) (other names
 15 include: Etilaam, Etizest, Depas, Etizola, Sedekopan, and
 16 Pasaden)
 17 Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-
 18 4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine)
 19 Gamma-hydroxybutyric acid (other names include GHB;
 20 gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium
 21 oxybate; sodium oxybutyrate) (2010)
 22 Mecloqualone (2572)
 23 Methaqualone (2565)
 24 (f) Stimulants. Unless specifically excepted or unless listed in
 25 another schedule, any material, compound, mixture, or preparation that
 26 contains any quantity of the following substances having a stimulant
 27 effect on the central nervous system, including its salts, isomers, and
 28 salts of isomers:
 29 ([+/-] cis-4-methylaminorex (([+/-])cis-4,5-
 30 dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)
 31 Aminorex (1585). Other names: aminoxaphen;
 32 2-amino-5-phenyl-2-oxazoline; or
 33 4,5-dihydro-5-phenyl-2-oxazolamine.
 34 Benzylone, 1-(1,3-benzodioxol-5-yl)-2-(benzylamino)propan-
 35 -1-one. Synonyms: BMDP, N-benzyl methylone,
 36 3,4-Methylene diox-Nbenzylcathinone,
 37 N-benzyl-3,4-methylenedioxcathinone.
 38 Cathinone (1235). Some trade or other names:
 39 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone;
 40 2-aminopropiophenone; and norephedrone.
 41 Fenethylamine (1503).
 42 N-Benzylpiperazine (7493). Other names: BZP; and



- 1 1-benzylpiperazine.
 2 N-ethylamphetamine (1475).
 3 Methcathinone (1237). Some other trade names:
 4 2-Methylamino-1-Phenylpropan-I-one; Ephedrone;
 5 Monomethylpropion; UR 1431.
 6 N, N-dimethylamphetamine (1480). Other names: N,
 7 N-alpha-trimethyl-benzeneethanamine; and N,
 8 N-alpha-trimethylphenethylamine.
 9 (g) Synthetic drugs as defined in IC 35-31.5-2-321.
 10 SECTION 8. IC 35-48-4-18 IS ADDED TO THE INDIANA CODE
 11 AS A NEW SECTION TO READ AS FOLLOWS [EFFECTIVE JULY
 12 1, 2022]: **Sec. 18. (a) This section applies to the following offenses:**
 13 **(1) Dealing in a schedule I, II, or III controlled substance**
 14 **under section 2 of this chapter.**
 15 **(2) Possession of a controlled substance under section 7 of this**
 16 **chapter.**
 17 **(3) Dealing in marijuana, hash oil, or hashish under section 10**
 18 **of this chapter.**
 19 **(4) Possession of marijuana, hash oil, or hashish under section**
 20 **11 of this chapter.**
 21 **(b) It is a defense to a prosecution for an offense described in**
 22 **subsection (a) that all of the following apply:**
 23 **(1) The controlled substance is hemp (as defined in 7 U.S.C.**
 24 **1639o(1)). (2) The controlled substance was being shipped**
 25 **from a licensed producer in another state in continuous**
 26 **transit through Indiana to a licensed handler in another state.**



COMMITTEE REPORT

Madam President: The Senate Committee on Corrections and Criminal Law, to which was referred Senate Bill No. 209, 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 35-31.5-2-321, AS AMENDED BY P.L.61-2020, SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 321. "Synthetic drug" means:

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

- (A) JWH-015 ((2-Methyl-1-propyl-1H-indol-3-yl)-1-naphthalenylmethanone).
- (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
- (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
- (D) JWH-073 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).
- (E) JWH-081 (4-methoxynaphthalen-1-yl-(1-pentylindol-3-yl)methanone).
- (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
- (G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-yl-methanone).
- (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
- (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
- (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
- (K) HU-210 ((6aR,10aR)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
- (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
- (M) HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-(2-methyloctan-2-yl)phenyl]-7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
- (N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-2,5-cyclohexadiene-1,4-dione).
- (O) CP 55,940 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]-5-(2-methyloctan-2-yl)phenol).



- (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol) and its homologues, or 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol), where side chain n=5, and homologues where side chain n=4, 6, or 7.
- (Q) WIN 55212-2
- ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-de)- 1,4- benzoxazin-6-yl]-1-naphthalenylmethanone).
- (R) RCS-4 ((4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone).
- (S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl)ethanone).
- (T) 4-Methylmethcathinone. Other name: mephedrone.
- (U) 3,4-Methylenedioxymethcathinone. Other name: methylone.
- (V) Fluoromethcathinone.
- (W) 4-Methoxymethcathinone. Other name: methedrone.
- (X) 4-Ethylmethcathinone (4-EMC).
- (Y) Methylenedioxypropylvalerone. Other name: MDPV.
- (Z) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
- (AA) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
- (BB) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
- (CC) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
- (DD) JWH-201, or 1-pentyl-3-(4-methoxyphenylacetyl)indole.
- (EE) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole.
- (FF) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
- (GG) CP 50,556-1, or [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpenta-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1-yl] acetate.
- (HH) Dimethylheptylpyran, or DMHP.
- (II) 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP.
- (JJ) 6-APB [6-(2-aminopropyl)benzofuran].
- (LL) 7-hydroxymitragynine.
- (MM) α -PPP [α -pyrrolidinopropiophenone].
- (NN) α -PVP (desmethylpyrovalerone).
- (OO) AM-251.



(PP) AM-1241.
 (QQ) AM-2201.
 (RR) AM-2233.
 (SS) Buphedrone (α -methylamino-butyrophenone (MABP)).
 (TT) Butylone.
 (UU) CP-47,497-C7.
 (VV) CP-47,497-C8.
 (WW) Desoxypipradol.
 (XX) Ethylone.
 (YY) Eutylone.
 (ZZ) Flephedrone.
 (AAA) JWH-011.
 (BBB) JWH-020.
 (CCC) JWH-022.
 (DDD) JWH-030.
 (EEE) JWH-182.
 (FFF) JWH-302.
 (GGG) MDAI [5,6-methylenedioxy-2-aminoindane].
 (HHH) Mitragynine.
 (III) Naphyrone.
 (JJJ) Pentedrone.
 (LLL) Pentylone.
 (MMM) Methoxetamine
 [2-(3-methoxyphenyl)-2-(ethylamino)-cyclohexanone].
 (NNN) A796,260 [1-(2-morpholin-4-ylethyl)-1H-indol-3-yl]-
 (2,2,3,3-tetramethylcyclopropyl)methanone].
 (OOO) AB-001[(1s,3s)-adamantan-1-yl]
 (1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
 (1-adamantoyl)indole].
 (PPP) AM-356 [Methanandamide].
 (QQQ) AM 1248 [1-[(1-methyl-2-piperidinyl) methyl]-
 1H-indol-3-yl] tricyclo[3.3.1.1^{3,7}] dec-1-yl-methanone]or
 [(1-[(N-methylpiperindin-2-yl)
 Methyl]-3-(Adamant-1-oyl)indole].
 (RRR) AM 2233 Azepane isomer [(2-iodophenyl)
 (1-(1-methylazepan-3-yl)-1H-indol-3-yl)methanone].
 (SSS) CB-13 [1-Naphthalenyl
 [4-(pentyoxy)-1-naphthalenyl]methanone].
 (TTT) UR-144 [(1-pentyl-1H-indol-3-yl)
 (2,2,3,3-tetramethylcyclopropyl)-methanone].
 (UUU) URB 597 [(3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl)-
 cyclohexylcarbamate].



(VVV) URB602 [[1,1'-biphenyl]-3-yl-carbamic acid, cyclohexyl ester].

(WWW) URB 754 [6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one].

(XXX) XLR-11 or 5-fluoro UR-144 (1-(5-fluoropentyl)-1H-indol-3-yl) (2,2,3,3-tetramethylcyclopropyl)methanone].

(YYY) AKB48 (Other names include: N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide; 1-pentyl-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-indazole-3-carboxamide).

(ZZZ) 25I-NBOMe (Other names include: 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine); 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine).

(AAAA) 2C-C-NBOMe (Other names include: 25C-NBOMe; 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 2,5-Dimethoxy-4-chloro-N-(2-methoxybenzyl)phenethylamine).

(BBBB) 2NE-1 (Other names include: 1-Pentyl-3-(1-adamantylamido)indole).

(CCCC) STS-135 (Other names include: N-Adamantyl-1-fluoropentylindole-3-carboxamide (1-5-fluoropentyl)-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-indole-3-carboxamide).

(DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinyl ester-1H-indole-2-carboxylic acid).

(EEEE) 5-Fluoro-PB-22 (Other names include: 1-(5-Fluoropentyl)-8-quinolinyl ester-1H-indole-3-carboxylic acid).

(FFFF) Benocyclidine (Other names include: BCP, BTCP, and Benzothiophenylcyclohexylpiperidine).

(GGGG) 25B-NBOMe (Other names include: 2C-B-NBOMe and 4-Bromo-2,5-dimethoxy-N-[(2-Methoxyphenyl)methyl]benzeneethanamine).

(HHHH) APB (Other names include: (2-Aminopropyl)Benzofuran).

(IIII) AB-PINACA (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-



indazole-3-carboxamide.
 (JJJJ) AB-FUBINACA
 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide).
 (KKKK) ADB-PINACA
 (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide).
 (LLLL) Fluoro ADBICA (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-(fluoropentyl)-1H-indole-3-carboxamide).
 (MMMM) APDB (Other names include: -EMA, -Desoxy-MDA, and (2-Aminopropyl)-2,3-dihydrobenzofuran).
 (NNNN) THJ-2201 (Other names include: AM2201 indazole analog, Fluoropentyl-JWH-018 indazole, and 5-Fluoro-THJ-018).
 (OOOO) AM 2201 benzimidazole analog (Other names include: FUBIMINA, FTHJ, and (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-yl)(naphthalene-1-yl)methanone).
 (PPPP) MN-25 (Other names include: 7-methoxy-1-[2-(4-morpholinyl)ethyl]-N-[1S, 2S, 4R)-1,3,3-trimethylbicyclo[2.2.1]hept-2-yl]-1H-indole-3-carboxamide and UR-12).
 (QQQQ) FUB-PB-22 (Other names include: Quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 (RRRR) FUD-PB-22 (Other names include: Naphthalen-1-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
 (SSSS) 5-Fluoro-AB-PINACA (Other names include: AB-PINACA 5-fluoro analog and N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide).
 (TTTT) 4-MePPP (Other names include: 4-methyl-alpha-pyrrolidinopropiophenone).
 (UUUU) alpha-PBP (Other names include: Alpha-pyrrolidinobutiophenone).
 (VVVV) AB-CHMINACA (Other names include: (N-[1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide).
 (WWWW) Mexedrone
 (3-methoxy-2-(methylamino)-1-(p-tolyl)propan-1-one).
 (XXXX) MT-45,
 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine).



- (YYYY) methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate [5F-ADB; 5F-MDMB-PINACA].
- (ZZZZ) methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate [5F-AMB].
- (AAAAA) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide [ADB-FUBINACA].
- (BBBBB) N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide [5F-APINACA, 5F-AKB48].
- (CCCCC) methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate [MDMB-CHMICA, MMB-CHMINACA].
- (DDDDD) methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate [MDMB-FUBINACA].
- (EEEEE) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide [MAB-CHMINACA and ADB-CHMINACA].
- (FFFFF) Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate [FUB-AMB, MMB-FUBINACA, AMB-FUBINACA].
- (GGGGG) 3,4-dichloro-N-[(1-dimethylamino)cyclohexylmethyl]benzamide [AH7921].
- (HHHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (trivial name: NM2201; CBL2201)
- (IIIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (trivial name: 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINACA; 4-CN-CUMYL BINACA; CUMYL-4CN-BINACA; SGT-78).
- (JJJJJ) methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoate (trivial names: MMB-CHMICA, AMB-CHMICA).
- (KKKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide (trivial name: 5F-CUMYL-P7AICA).
- (LLLLL) N-1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (N-ethylpentylone, ephylone).
- (MMMMM) Synthetic cathinone, 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one (N-ethylpentylone, ephylone) and its optical, positional, and geometric isomers, salts, and salts of isomers.



(NNNNN) ethyl
 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (trivial name: 5F-EDMB-PINACA).
 (OOOOO) methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate (trivial name: 5F-MDMB-PICA).
 (PPPPP) N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (trivial names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-FLUOROBENZYL)).
 (QQQQQ) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (trivial names: 5F-CUMYL-PINACA; SGT-25).
 (RRRRR) 1-(4-fluorobenzyl)-1H-indol-3-yl(2,2,3,3-tetramethylcyclopropyl) methanone (trivial name: FUB-144).
 (SSSSS) 4F-MDMB-BINACA.
 (TTTTT) N-ethylhexedrone
 (2-(ethylamino)-1-phenylhexan-1-one).
 (UUUUU) alpha-pyrrolidinohexanophenone
 (1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one).
 (VVVVV) alpha-pyrrolidinohexiophenone; trivial name: a-PHP.
 (WWWWW) 4'-methyl-alpha-pyrrolidinohexiophenone
 (1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one).
 (XXXXX) 4-methyl-alphaethylaminopentiophenone
 (2-(ethylamino)-1-(4-methylphenyl)pentan-1-one; trivial name: 4-MEAP.
 (YYYYY) 4'-methyl-alpha-pyrrolidinohexanophenone; trivial name: MPHP.
 (ZZZZZ) alpha-pyrrolidinoheptaphenone
 (1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one; trivial name: PV8.
 (AAAAA) 4'-chloro-alpha-pyrrolidinovalerophenone (1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one).
 (BBBBB) 4'-chloro-alpha-pyrrolidinopentiophenone; trivial name: 4-chloro-a-PVP.
(CCCCCC) 4,4'-dimethylaminorex (common name: 4,4'-DMAR).

(2) Any compound structurally derived from 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidiny)methyl, 2-(4-morpholinyl)ethyl, or



1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent.

(3) Any compound structurally derived from 3-(1-naphthoyl)pyrrole by substitution at the nitrogen atom of the pyrrole ring by alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the pyrrole ring to any extent and whether or not substituted in the naphthyl ring to any extent.

(4) Any compound structurally derived from 1-(1-naphthylmethyl)indene by substitution at the 3-position of the indene ring by alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent.

(5) Any compound structurally derived from 3-phenylacetylindole by substitution at the nitrogen atom of the indole ring with alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent.

(6) Any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not substituted in the cyclohexyl ring to any extent.

(7) Any compound containing a 3-(benzoyl)indole structure with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or



1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent.

(8) Any compound, except bupropion or a compound listed under a different schedule, structurally derived from 2-aminopropan-1-one by substitution at the 1-position with either phenyl, naphthyl, or thiophene ring systems, whether or not the compound is further modified:

(A) by substitution in the ring system to any extent with alkyl, alkylendioxy, alkoxy, haloalkyl, hydroxyl, or halide substituents, whether or not further substituted in the ring system by one or more other univalent substituents;

(B) by substitution at the 3-position with an acyclic alkyl substituent;

(C) by substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or methoxybenzyl groups; or

(D) by inclusion of the 2-amino nitrogen atom in a cyclic structure.

(9) Any compound structurally derived from 3-tetramethyl cyclopropanoylindole with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the tetramethylcyclopropyl ring to any extent.

(10) Any compound containing a N-(1-adamantyl)-1H-indazole-3-carboxamide structure with substitution at the nitrogen atom of the indazole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted at the nitrogen atom of the carboxamide to any extent, whether or not further substituted in the indazole ring to any extent, and whether or not further substituted on the adamantyl ring system to any extent. An example of this structural class includes AKB48.

(11) Any compound containing a N-(1-adamantyl)-



1H-indole-3-carboxamide structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted at the nitrogen atom of the carboxamide to any extent, whether or not further substituted in the indole ring to any extent, and whether or not further substituted on the adamantyl ring system to any extent. An example of this structural class includes STS-135.

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

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

SECTION 2. IC 35-48-1-17.4, AS ADDED BY P.L.61-2020, SECTION 3, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 17.4. (a) Except as provided in subsections (b) and (c), "isomer" means an optical isomer.

(b) "Isomer", as used in IC 35-48-2-4(d), means an optical, positional, or geometric isomer.

(c) "Isomer", as used in section 7 of this chapter, means an optical or geometric isomer. "Isomer" means an optical, positional, or geometric isomer."

Page 2, between lines 17 and 18, begin a new line block indented and insert:

"Brorphine".

Page 2, delete line 42, begin a new line block indented, and insert:

"2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-benzimidazole (N-pyrrolidino etonitazene; etonitazepyne)".

Page 3, delete line 1.

Page 6, strike line 1.



Page 6, line 2, strike "includes the optical, position, and geometric isomers):" and insert "designation:".

Re-number all SECTIONS consecutively.

and when so amended that said bill do pass.

(Reference is to SB 209 as introduced.)

YOUNG M, Chairperson

Committee Vote: Yeas 8, Nays 0.

SENATE MOTION

Madam President: I move that Senate Bill 209 be amended to read as follows:

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

"SECTION 1. IC 15-15-13-6.5, AS ADDED BY P.L.190-2019, SECTION 7, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 6.5. As used in this chapter, "hemp product" means a product derived from, or made by, processing hemp plants or plant parts including derivatives, extracts, cannabinoids, isomers, acids, salts, and salts of isomers. However, the term does not include:

- (1) smokable hemp (as defined by IC 35-48-1-26.6); or
- (2) products that contain a total ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** concentration of more than three-tenths of one percent (0.3%) by weight.

SECTION 2. IC 24-4-21-3, AS AMENDED BY P.L.190-2019, SECTION 24, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 3. A person may distribute low THC hemp extract in Indiana only if the distributor has a certificate of analysis prepared by an independent testing laboratory showing:

- (1) that the low THC hemp extract is the product of a batch tested by the independent testing laboratory;
- (2) that the independent testing laboratory determined that the batch contained not more than three-tenths percent (0.3%) total ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** (THC), including precursors **and isomers**, by weight, based on the testing of a random sample of the batch; and
- (3) the cannabidiol percent present of the low THC hemp extract.

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SECTION 3. IC 24-4-21-4, AS ADDED BY P.L.153-2018, SECTION 9, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 4. (a) Except as provided in subsection (b), low THC hemp extract must be distributed in packaging that contains the following information:

(1) A scannable bar code or QR code linked to a document that contains information with respect to the manufacture of the low THC hemp extract, including the:

- (A) batch identification number;
- (B) product name;
- (C) batch date;
- (D) expiration date, which must be not more than two (2) years from the date of manufacture;
- (E) batch size;
- (F) total quantity produced;
- (G) ingredients used, including the:
 - (i) ingredient name;
 - (ii) name of the company that manufactured the ingredient;
 - (iii) company or product identification number or code, if applicable; and
 - (iv) ingredient lot number; and
- (H) download link for a certificate of analysis for the low THC hemp extract.

(2) The batch number.

(3) The Internet address of a web site to obtain batch information.

(4) The expiration date.

(5) The number of milligrams of low THC hemp extract.

(6) The manufacturer.

(7) The fact that the product contains not more than three-tenths percent (0.3%) total ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** (THC), including precursors **and isomers**, by weight.

(b) Before July 1, 2018, low THC hemp extract may be distributed in Indiana without having met the requirements described in subsection (a)."

Page 1, line 5, delete "compound:" and insert "compound **and its isomers, salts, and salts of isomers:**".

Page 11, between lines 5 and 6, begin a new paragraph and insert:

"SECTION 6. IC 35-48-1-17.5, AS AMENDED BY P.L.190-2019, SECTION 28, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: Sec. 17.5. (a) "Low THC hemp extract" means a substance or compound that:



- (1) is derived from or contains any part of the plant *Cannabis sativa* L. that meets the definition of hemp under IC 15-15-13-6;
 - (2) contains not more than three-tenths percent (0.3%) total ~~delta-9-tetrahydrocannabinol~~ **tetrahydrocannabinol** (THC), including precursors **and isomers**, by weight; and
 - (3) contains no other controlled substances.
- (b) The term does not include:
- (1) the harvested reproductive organ, whether immature or mature, of the female hemp plant; or
 - (2) smokable hemp."
- Renumber all SECTIONS consecutively.
- (Reference is to SB 209 as printed January 21, 2022.)

BROWN L

SENATE MOTION

Madam President: I move that Engrossed Senate Bill 209, which is eligible for third reading, be returned to second reading for purposes of amendment.

YOUNG M

SENATE MOTION

Madam President: I move that Senate Bill 209 be amended to read as follows:

Page 21, after line 9, begin a new paragraph and insert:

"SECTION 8. IC 35-48-4-18 IS ADDED TO THE INDIANA CODE AS A NEW SECTION TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: **Sec. 18. (a) This section applies to the following offenses:**

- (1) Dealing in a schedule I, II, or III controlled substance under section 2 of this chapter.**
- (2) Possession of a controlled substance under section 7 of this chapter.**
- (3) Dealing in marijuana, hash oil, or hashish under section 10 of this chapter.**
- (4) Possession of marijuana, hash oil, or hashish under section**



11 of this chapter.

(b) It is a defense to a prosecution for an offense described in subsection (a) that all of the following apply:

(1) The controlled substance is hemp (as defined in 7 U.S.C. 1639o(1)). (2) The controlled substance was being shipped from a licensed producer in another state in continuous transit through Indiana to a licensed handler in another state."

Renumber all SECTIONS consecutively.

(Reference is to SB 209 as reprinted January 25, 2022.)

YOUNG M

