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.

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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:

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.

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).

SECTION 4. 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 **and its isomers, salts, and salts of isomers:**

(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-napthalenylmethanone).

(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) Methylenedioxypyrovalerone. 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-phenylpe
ntan-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)-admantan-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[3.3.1.13,7] 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[3.3.1.13.7]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[3.3.1.13.7]dec-1-yl-1H-indole-3-carboxamide).

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

(EEEE) 5-Fluoro-PB-22 (Other names include: 1-(5-Fluropentyl)-8-quinolinyl ester1H-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-dimenthoxy-N-[(2-Methozyphenyl)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-indaole-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-methyl1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxaminde).
(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-[(1dimethylamino)cyclohexylmethyl]benzamide) [AH7921].

(HHHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1 H-indole-3-carboxylate (trivial name: NM2201; CBL2201)

(IIIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1 H-indazole-3-carboxamide (trivial name: 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINA CA; 4-CN-CUMYL BINACA; CUMYL-4CN-BINACA; SGT-78).

(JJJJJ) methyl 2-(1-(cyclohexylmethyl)-1 H-indole-3-carboxamido)-3-methylbutanoate (trivial names: MMB-CHMICA, AMB-CHMICA).

(KKKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1 H-pyrrolo[2,3-b]pyridine-3-carboxamide (trivial name: 5F-CUMYL-P7AICA).

(LLLLL) N-1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-p entanone (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-alphapyrrolidinohexanophenone; trivial name: MPHP.

(ZZZZZ) alphapyrrolidinoheptaphenone (1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one; trivial name: PV8.

(AAAAAA) 4'-chloro-alphapyrrolidinovalerophenone (1-(4-chlorophenyl)-2-(pyrrolidin-1- yl)pentan-1-one.

(BBBBBB) 4'-chloro-alphapyrrolidinopentiophenone; 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-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 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, alkylenedioxy, 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 5. 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.**

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.

SECTION 7. IC 35-48-2-4, AS AMENDED BY P.L.10-2021, SECTION 1, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2022]: 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)

Brorphine

2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1yl)-N,N-diethylethan-1-amine (butonitazene)

Clonitazene (9612)

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

Dextromoramide (9613)

Diampromide (9615)

Diethylthiambutene (9616)

N,N-diethyl-2-(2-(4-flourobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine (flunitazene)

N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine (metodesnitazene)

N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidaz

ol-1-yl)ethan-1-amine (metonitazene)
N,N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1H-benzimidaz
ol-1-yl)ethan-1-amine (protonitazene)
Difenoxin (9168)
Dimenoxadol (9617)
Dimepheptanol (9618)
Dimethylthiambutene (9619)
Dioxaphetyl butyrate (9621)
Dipipanone (9622)
2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1yl)-N,N-diethyleth
an-1-amine (etodesnitazene; etazene)
2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-
benzimidazole (N-pyrrolidino etonitazene; etonitazepyne)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxeridine (9625)
Fentanyl related substances.
Furanyl fentanyl.
Furethidine (9626)
Hydroxypethidine (9627)
Isobutyryl fentanyl. Other name: N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide
Isotonitazene. Other name: N,N-diethyl-2-(2-(4 isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine)
Ketobemidone (9628)
Levomoramide (9629)
Levophenacylmorphan (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-propanimide](9813)
3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9833)
MDMB-4en-PINACA
4F-MDMB-BICA; 4-fluoro MDMB-BICA; 4F-MDMB-BUTICA; Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate
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-phenethylpiperidin-4-yl)- Nphenylcyclopentanecarboxamide (cyclopentyl fentanyl)

Noracymethadol (9633)

Norlevorphanol (9634)

Normethadone (9635)

Norpipanone (9636)

Ocfentanil. Other name: 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-carboxamide
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)
Valeryl fentanyl. Other name: 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)
Hydromorphinol (9301)
Methyldesorphine (9302)
Methyldihydromorphine (9304)
Morphine methylbromide (9305)
Morphine methylsulfonate (9306)
Morphine-N-Oxide (9307)

Myrophine (9308)
Nicocodeine (9309)
Nicomorphine (9312)
Normorphine (9313)
Pholcodine (9314)
Thebacon (9315)

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

(1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name: TCPy.

(2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or other names: 4-Bromo-2, 5-Dimethoxy-a-methylphenethylamine; 4-Bromo-2, 5-DMA.

(3) 4-Bromo-2, 5-dimethoxyphenethylamine (7392). Some trade or other names:
2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B, Nexus.

(4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name: DOET.

(5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348). Other name: 2C-T-7.

(6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.

(7) 4-Methoxyamphetamine (7411). Some trade or other names: 4-Methoxy-a-methylphenethylamine; Paramethoxyamphetamine; PMA.

(8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401). Other Name: MMDA.

(9) 5-Methoxy-N, N-diisopropyltryptamine, including any isomers, salts, or salts of isomers (7439). Other name: 5-MeO-DIPT.

(10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade and other names: 4-methyl-2,
5-dimethoxy-a-methylphenethylamine; DOM; and STP.

(11) 3, 4-methylenedioxy amphetamine (7400). Other name: MDA.

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

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

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

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

(C) $\pi^{3,4}$ 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] [1,4diazepine) (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.

Benzylone, 1-(1,3-benzodioxol-5-yl)-2-(benzylamino)propan-1-one. Synonyms: BMDP, N-benzyl methylone, 3,4-Methylenedioxy-Nbenzylcathinone, N-benzyl-3,4-methylenedioxycathinone.

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

Fenethylline (1503).

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

1-benzylpiperazine.

N-ethylamphetamine (1475).

Methcathinone (1237). Some other trade names: 2-Methylamino-1-Phenylpropan-I-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.

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.

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-napthalenylmethanone).
(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) Methylenedioxypyrovalerone. 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-phenylpe
ntan-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)-admantan-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[3.3.1.13,7] 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[3.3.1.13.7]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[3.3.1.13.7]dec-1-yl-1H-indole-3-carboxamide).
(DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinlyl ester-1H-indole-2-carboxylic acid).
(EEEE) 5-Fluoro-PB-22 (Other names include: 1-(5-Fluropentyl)-8-quinolinyl ester1H-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-dimenthoxy-N-[(2-Methozyphenyl)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-indaole-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-methyl1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxaminde).
(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-[(1dimethylamino)cyclohexylmethyl]benzamide) [AH7921].
(HHHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1 H-indole-3-carboxylate (trivial name: NM2201; CBL2201)
(IIIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1 H-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)-1 H-indole-3-carboxamido)-3-methylbutanoate (trivial names: MMB-CHMICA, AMB-CHMICA).
(KKKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1 H-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-alphapyrrolidinohexanophenone; trivial name: MPHP.

(ZZZZZ) alphapyrrolidinoheptaphenone (1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one; trivial name: PV8.

(AAAAAA) 4'-chloro-alphapyrrolidinovalerophenone (1-(4-chlorophenyl)-2-(pyrrolidin-1- yl)pentan-1-one.

(BBBBBB) 4'-chloro-alphapyrrolidinopentiophenone; 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-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 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, alkylenedioxy, 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:".

Renumber 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.

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