SENATE BILL No. 74

DIGEST OF SB 74 (Updated January 9, 2018 10:59 am - DI 84)

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

Synopsis: Controlled substances. Adds the substance Mexedrone to the definition of "synthetic drug" and adds additional controlled substances to the existing statutory list of depressants, hallucinogens, and opiates classified as schedule I.

Effective: July 1, 2018.

Young M

January 3, 2018, read first time and referred to Committee on Corrections and Criminal Law.
January 9, 2018, reported favorably — Do Pass.

SB 74—LS 6233/DI 131
SENATE BILL No. 74

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 35-31.5-2-321, AS AMENDED BY P.L.8-2017, SECTION 1, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2018]: 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-yethyl)indol-3-yl)-naphthalen-1-yl-methanone).
(H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetil)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) Methylenedioxyxypovalerone. 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-naphthyl)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-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1-yl] acetate.

(HH) Dimethylheptylpyran, or DMHP.

(II) 4-Methyl-alpha-pyrrolidinobutylphenone, 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) Desoxyxipipradol.

(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) Pentyldone.

(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].
(1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
(1-adamantoyl)indole].
(Methanandamide].
[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]).
[2-(iodophenyl)
1H-indol-3-yl)methanone].
[4-(pentoxy)- 1-naphthalenyl]methanone].
[(1-pentyl-1H-indol-3-yl)]
(2,2,3,3-tetramethylcyclopropyl)-methanone].
[3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl]-
cyclohexylcarbamate].
[[1,1'-biphenyl]- 3-yl-carbamic acid,
cyclohexyl ester].
[6-methyl-2-[(4-methylphenyl)
amino]-1-benzoazin-4-one].
5-fluoro UR-144
(1-(5-fluoropentyl)-1H-indol-3-yl)
[(2,2,3,3-tetramethylcyclopropyl)methanone].
N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide;
1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indazole-3-
carboxamide).
251-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).
(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).
(2NE-1 (Other names include: 1-Pentyl-3-
(1-adamantylamido)indole).
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-quinolinyl ester-1H-indole-2-carboxylic acid).

(EEEE) 5-Fluoro-PB-22 (Other names include: 1-(5-Fluoropentyl)-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-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, FTHI, 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-morpholiny1)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-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).
(DDDD) Acetyl fentanyl (Other names include:
N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide).
(XXXX) Mexedrone (3-methoxy-2-(methylamino)-1-(p-tolyl)propan-1-one).

(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, 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, 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 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, 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 35-48-2-4, AS AMENDED BY P.L.8-2017, SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2018]: Sec. 4. (a) The controlled substances listed in this section are included in schedule I.

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

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

Acetylmethadol (9601)

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

Allylprodine (9602)

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

Alphacetylmethadol (9603)

Alphameprodine (9604)

Alphamethadol (9605)

Alphamethylfentanyl (9814)

Benzethidine (9606)

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

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

Betacetylmethadol (9607)

Betameprodine (9608)

Betamethadol (9609)

Betaprodine (9611)

Clonitazene (9612)

Dextromoramide (9613)

Diampromide (9615)

Diethylthiambutene (9616)

Difenoxin (9168)

Dimenoxadol (9617)

Dimephtanol (9618)

Dimethylthiambutene (9619)

Dioxaphetyl butyrate (9621)
Dipipanone (9622)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxeridine (9625)
Furethidine (9626)
Hydroxypethidine (9627)
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)
MPPP (1-methyl-4-phenyl-4-propionoxy Piperidine) (9961)

Morpheridine (9632)
N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), including any isomers, salts, or salts of isomers (9818)
N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), including any isomers, salts, or salts of isomers (9834)

Noracymethadol (9633)
Norlevorphanol (9634)
Normethadone (9635)
Norpipanone (9636)

**Ortho-fluorofentanyl or 2-fluorofentanyl. Other name:**
N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide
Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide) (9812)
Phenadoxone (9637)
Phenampromide (9638)
Phenomorphan (9647)
Phenoperidine (9641)
PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxy piperidine] (9663)
Piritramide (9642)
Proheptazine (9643)
Properidine (9644)
Propiram (9649)
Racemoramide (9645)

**Tetrahydrofuranyl fentanyl. Other name:**

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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)

(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:

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

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

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

(2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or
other names: 4-Bromo-2, 5-Dimethoxy-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-ethylamphetamine (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-methylenedioxyamphetamine (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-methylenedioxyamphetamine (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-trimethoxyamphetamine (7390). Other name: TMA.

(15) Alpha-ethyltryptamine (7249). Some trade and other names:
Etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine;
3-(2-aminoethyl) 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 cis or trans tetrahydrocannabinol, and their optical isomers;
   (B) π cis or trans tetrahydrocannabinol, and their optical isomers; and
   (C) π 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-phenylethylcyclohexylamine; (1-phenylethylcyclohexyl) ethylamine; N-(1-phenylethylcyclohexyl) ethylamine; cyclohexamine; PCE.

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

(34) Thiophene analog of phencyclidine (7470). Some trade or other names: 1-(1-(2-thienyl)cyclohexyl) piperidine; 2-Thienyl Analogue 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 SB 74—LS 6233/DI 131
specific chemical designation:

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

Flubromazolam (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine)

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

Mecloqualone (2572)

Methaqualone (2565)

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

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

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

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

Fenethylline (1503).

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

N-ethylamphetamine (1475).

Methcathinone (1237). Some other trade names: 2-Methylamino-1-Phenylpropan-1-one; Ephedrine; 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.
COMMITTEE REPORT

Madam President: The Senate Committee on Corrections and
Criminal Law, to which was referred Senate Bill No. 74, has had the
same under consideration and begs leave to report the same back to the
Senate with the recommendation that said bill DO PASS.

(Reference is to SB 74 as introduced.)

YOUNG M, Chairperson

Committee Vote: Yeas 8, Nays 0