March 15, 2017

ENGROSSED
HOUSE BILL No. 1019

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

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

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

Effective: July 1, 2017.

Ellington, Negele, Klinker, Hatfield

(SENATE SPONSORS — MERRITT, HEAD, GROOMS)

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

SENATE ACTION

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

March 15, 2017

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

PRINTING CODE. Amendments: Whenever an existing statute (or a section of the Indiana Constitution) is being amended, the text of the existing provision will appear in this style type, additions will appear in **this style type**, and deletions will appear in ~~this style type.~~

Additions: Whenever a new statutory provision is being enacted (or a new constitutional provision adopted), the text of the new provision will appear in **this style type**. Also, the word **NEW** will appear in that style type in the introductory clause of each SECTION that adds a new provision to the Indiana Code or the Indiana Constitution.

Conflict reconciliation: Text in a statute in *this style type* or ~~*this style type*~~ reconciles conflicts between statutes enacted by the 2016 Regular Session of the General Assembly.

ENGROSSED
HOUSE BILL No. 1019

A BILL FOR AN ACT to amend the Indiana Code concerning criminal law and procedure.

Be it enacted by the General Assembly of the State of Indiana:

SECTION 1. IC 35-31.5-2-321, AS AMENDED BY P.L.168-2016, SECTION 10, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2017]: 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) Acetyl fentanyl (Other names include: N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide).

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

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

Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide) (9815)
Acetylmethadol (9601)
Allylprodine (9602)
Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)
Alphacetylmethadol (9603)
Alphameprodine (9604)
Alphamethadol (9605)
Alphamethylfentanyl (9814)
Benzethidine (9606)
Beta-hydroxy-3-methylfentanyl (9831). Other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide
Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)
Betacetylmethadol (9607)
Betameprodine (9608)
Betamethadol (9609)
Betaprodine (9611)
Clonitazene (9612)
Dextromoramide (9613)
Diampromide (9615)
Diethylthiambutene (9616)
Difenoxin (9168)
Dimenoxadol (9617)
Dimepheptanol (9618)
Dimethylthiambutene (9619)
Dioxaphetyl butyrate (9621)
Dipipanone (9622)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxeridine (9625)

Furethidine (9626)
Hydroxypethidine (9627)
Ketobemidone (9628)
Levomoramide (9629)
Levophenacylmorphan (9631)
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-propionoxypiperidine) (9961)
Morpheridine (9632)
N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), including any isomers, salts, or salts of isomers (9818)
N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), including any isomers, salts, or salts of isomers (9834)
Noracymethadol (9633)
Norlevorphanol (9634)
Normethadone (9635)
Norpipanone (9636)
Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
Phenadoxone (9637)
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)
Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide) (9835)
Tilidine (9750)
Trimeperidine (9646)
U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-N-methyl- benzamide)

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

the specific chemical designation:

Acetorphine (9319)
Acetyldihydrocodeine (9051)
Benzylmorphine (9052)
Codeine methylbromide (9070)
Codeine-N-Oxide (9053)
Cyprenorphine (9054)
Desomorphine (9055)
Dihydromorphine (9145)
Drotebanol (9335)
Etorphine (except hydrochloride salt) (9056)
Heroin (9200)
Hydromorphinol (9301)
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):

(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).

(e) Depressants. Unless specifically excepted in a rule adopted by the board or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

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

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

Mecloqualone (2572)

Methaqualone (2565)

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

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

Aminorex (1585). Other names: aminoxaphen;

2-amino-5-phenyl-2-oxazoline; or
4,5-dihydro-5-phenyl-2-oxazolamine.
Cathinone (1235). Some trade or other names: 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone; 2-aminopropiophenone; and norephedrone.
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.

COMMITTEE REPORT

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

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

Page 6, delete line 16.

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

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

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

Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide) (9815)
Acetylmethadol (9601)
Allylprodine (9602)
Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)
Alphacetylmethadol (9603)
Alphameprodine (9604)
Alphamethadol (9605)
Alphamethylfentanyl (9814)
Benzethidine (9606)
Beta-hydroxy-3-methylfentanyl (9831). Other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide
Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)
Betacetylmethadol (9607)
Betameprodine (9608)
Betamethadol (9609)
Betaprodine (9611)
Clonitazene (9612)
Dextromoramide (9613)
Diampromide (9615)

Diethylthiambutene (9616)
Difenoxin (9168)
Dimenoxadol (9617)
Dimepheptanol (9618)
Dimethylthiambutene (9619)
Dioxaphetyl butyrate (9621)
Dipipanone (9622)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxeridine (9625)
Furethidine (9626)
Hydroxypethidine (9627)
Ketobemidone (9628)
Levomoramide (9629)
Levophenacylmorphan (9631)
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-propionoxypiperidine) (9961)
Morpheridine (9632)
N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), including any isomers, salts, or salts of isomers (9818)
N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), including any isomers, salts, or salts of isomers (9834)
Noracymethadol (9633)
Norlevorphanol (9634)
Normethadone (9635)
Norpipanone (9636)
Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
Phenadoxone (9637)
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)

Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide) (9835)
Tilidine (9750)
Trimeperidine (9646)
U47700 (3,4-dichloro- N- [2-dimethylamino)cyclohexyl]-N-methyl- benzamide)

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

Acetorphine (9319)
Acetyldihydrocodeine (9051)
Benzylmorphine (9052)
Codeine methylbromide (9070)
Codeine-N-Oxide (9053)
Cyprenorphine (9054)
Desomorphine (9055)
Dihydromorphine (9145)
Drotebanol (9335)
Etorphine (except hydrochloride salt) (9056)
Heroin (9200)
Hydromorphinol (9301)
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):

(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).

(e) Depressants. Unless specifically excepted in a rule adopted by the board or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

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

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

Mecloqualone (2572)
Methaqualone (2565)

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

([+/-]) cis-4-methylaminorex (([+/-])cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)
Aminorex (1585). Other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine.
Cathinone (1235). Some trade or other names: 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone; 2-aminopropiophenone; and norephedrone.
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.".

Renumber all SECTIONS consecutively.

and when so amended that said bill do pass.

(Reference is to HB 1019 as introduced.)

WASHBURNE

Committee Vote: yeas 8, nays 0.

COMMITTEE REPORT

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

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

YOUNG M, Chairperson

Committee Vote: Yeas 7, Nays 0