

PROPOSED AMENDMENT

SB 209 # 1

DIGEST

Updates the scheduled drug list.

- 1 Page 1, between the enacting clause and line 1, begin a new
2 paragraph and insert:
3 "SECTION 1. IC 35-31.5-2-321, AS AMENDED BY P.L.61-2020,
4 SECTION 2, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE
5 JULY 1, 2022]: Sec. 321. "Synthetic drug" means:
6 (1) a substance containing one (1) or more of the following
7 chemical compounds, including an analog of the compound:
8 (A) JWH-015 ((2-Methyl-1-propyl-1H-
9 indol-3-yl)-1-naphthalenylmethanone).
10 (B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).
11 (C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).
12 (D) JWH-073 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).
13 (E) JWH-081 (4-methoxynaphthalen-1-yl- (1-pentylindol-
14 3-yl)methanone).
15 (F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).
16 (G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-
17 naphthalen-1-yl-methanone).
18 (H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
19 (I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
20 (J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
21 (K) HU-210 ((6aR,10aR)- 9-(Hydroxymethyl)- 6,6-dimethyl-
22 3-(2-methyloctan-2-yl)-
23 6a,7,10,10a-tetrahydrobenzo [c]chromen- 1-ol).
24 (L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)- 6,6-dimethyl-
25 3-(2-methyloctan-2-yl)- 6a,7,10,10a-tetrahydrobenzo
26 [c]chromen-1-ol).
27 (M) HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-
28 (2-methyloctan- 2-yl)phenyl]-

- 1 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
- 2 (N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-
- 3 (1-methylethenyl)-2 -cyclohexen-1-yl]-5
- 4 -pentyl-2,5-cyclohexadiene-1,4-dione).
- 5 (O) CP 55,940
- 6 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]-
- 7 5-(2-methyloctan-2-yl)phenol).
- 8 (P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]- 5-
- 9 (2-methyloctan-2-yl)phenol) and its homologues, or
- 10 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)
- 11 phenol), where side chain n=5, and homologues where side
- 12 chain n=4, 6, or 7.
- 13 (Q) WIN 55212-2
- 14 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
- 15 pyrrolo [1,2,3-de)- 1,4- benzoxazin-
- 16 6-yl]-1-naphthalenylmethanone).
- 17 (R) RCS-4 ((4-methoxyphenyl)
- 18 (1-pentyl-1H-indol-3-yl)methanone).
- 19 (S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-
- 20 indol-3-yl)-2-(2-methoxyphenyl)ethanone).
- 21 (T) 4-Methylmethcathinone. Other name: mephedrone.
- 22 (U) 3,4-Methylenedioxymethcathinone. Other name:
- 23 methylone.
- 24 (V) Fluoromethcathinone.
- 25 (W) 4-Methoxymethcathinone. Other name: methedrone.
- 26 (X) 4-Ethylmethcathinone (4-EMC).
- 27 (Y) Methylenedioxyprovalerone. Other name: MDPV.
- 28 (Z) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole.
- 29 (AA) JWH-098, or
- 30 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
- 31 (BB) JWH-164, or
- 32 1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
- 33 (CC) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
- 34 (DD) JWH-201, or
- 35 1-pentyl-3-(4-methoxyphenylacetyl)indole.
- 36 (EE) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole.
- 37 (FF) AM-694, or
- 38 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
- 39 (GG) CP 50,556-1, or
- 40 [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpe

- 1 ntan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1
 2 -yl] acetate.
 3 (HH) Dimethylheptylpyran, or DMHP.
 4 (II) 4-Methyl- α -pyrrolidinobutiophenone, or MPBP.
 5 (JJ) 6-APB [6-(2-aminopropyl)benzofuran].
 6 (LL) 7-hydroxymitragynine.
 7 (MM) α -PPP [α -pyrrolidinopropiophenone].
 8 (NN) α -PVP (desmethylpyrovalerone).
 9 (OO) AM-251.
 10 (PP) AM-1241.
 11 (QQ) AM-2201.
 12 (RR) AM-2233.
 13 (SS) Buphedrone (α -methylamino-butyrophenone (MABP)).
 14 (TT) Butylone.
 15 (UU) CP-47,497-C7.
 16 (VV) CP-47,497-C8.
 17 (WW) Desoxypipradol.
 18 (XX) Ethylone.
 19 (YY) Eutylone.
 20 (ZZ) Flephedrone.
 21 (AAA) JWH-011.
 22 (BBB) JWH-020.
 23 (CCC) JWH-022.
 24 (DDD) JWH-030.
 25 (EEE) JWH-182.
 26 (FFF) JWH-302.
 27 (GGG) MDAI [5,6-methylenedioxy-2-aminoindane].
 28 (HHH) Mitragynine.
 29 (III) Naphyrone.
 30 (JJJ) Pentedrone.
 31 (LLL) Pentylone.
 32 (MMM) Methoxetamine
 33 [2-(3-methoxyphenyl)-2-(ethylamino)-cyclohexanone].
 34 (NNN) A796,260 [1-(2-morpholin-4-ylethyl)-1H-indol-3-yl]-
 35 (2,2,3,3-tetramethylcyclopropyl)methanone].
 36 (OOO) AB-001[(1s,3s)-adamantan-1-yl]
 37 (1-pentyl-1H-indol-3-yl)methanone] or [1-Pentyl-3-
 38 (1-adamantoyl)indole].
 39 (PPP) AM-356 [Methanandamide].
 40 (QQQ) AM 1248 [1-[(1-methyl-2-piperidinyl) methyl]-

1 1H-indol-3-yl] tricyclo[3.3.1.1^{3,7}] dec-1-yl-methanone]or
 2 [(1-[(N-methylpiperindin-2-yl)
 3 Methyl]-3-(Adamant-1-oyl)indole].
 4 (RRR) AM 2233 Azepane isomer [(2-iodophenyl)
 5 (1-(1-methylazepan-3-yl)- 1H-indol-3-yl)methanone].
 6 (SSS) CB-13 [1-Naphthalenyl
 7 [4-(pentyoxy)- 1-naphthalenyl]methanone].
 8 (TTT) UR-144 [(1-pentyl-1H-indol-3-yl)
 9 (2,2,3,3-tetramethylcyclopropyl)-methanone].
 10 (UUU) URB 597 [(3'-(aminocarbonyl) [1,1'-biphenyl]-3-yl)-
 11 cyclohexylcarbamate].
 12 (VVV) URB602 [[1,1'-biphenyl]- 3-yl-carbamic acid,
 13 cyclohexyl ester].
 14 (WWW) URB 754 [6-methyl-2-[(4-methylphenyl)
 15 amino]-1-benzoxazin-4-one].
 16 (XXX) XLR-11 or 5-fluoro UR-144
 17 (1-(5-fluoropentyl)-1H-indol-3-yl)
 18 (2,2,3,3-tetramethylcyclopropyl)methanone].
 19 (YYY) AKB48 (Other names include:
 20 N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide;
 21 1-pentyl-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-indazole-3-
 22 carboxamide).
 23 (ZZZ) 25I-NBOMe (Other names include:
 24 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-
 25 benzeneethanamine);
 26 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
 27 methyl]ethanamine).
 28 (AAAA) 2C-C-NBOMe (Other names include: 25C-NBOMe;
 29 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
 30 methyl]ethanamine;
 31 2,5-Dimethoxy-4-chloro-N-(2-methoxybenzyl)
 32 phenethylamine).
 33 (BBBB) 2NE-1 (Other names include: 1-Pentyl-3-
 34 (1-adamantylamido)indole).
 35 (CCCC) STS-135 (Other names include:
 36 N-Adamantyl-1-fluoropentylindole-3- carboxamide
 37 (1-5-fluoropentyl)-N-tricyclo[3.3.1.1^{3,7}]dec-1-yl-1H-
 38 indole-3-carboxamide).
 39 (DDDD) PB-22 (Other names include: 1-Pentyl-8-quinolinyl
 40 ester-1H-indole-2-carboxylic acid).

1 (EEEE) 5-Fluoro-PB-22 (Other names include:
2 1-(5-Fluoropentyl)-8-quinolinyl ester 1H-indole-3-carboxylic
3 acid).
4 (FFFF) Benocyclidine (Other names include: BCP, BTCP, and
5 Benzothiophenylcyclohexylpiperidine).
6 (GGGG) 25B-NBOMe (Other names include: 2C-B-NBOMe
7 and 4-Bromo-2,
8 5-dimethoxy-N-[(2-Methoxyphenyl)methyl]
9 benzenethanamine).
10 (HHHH) APB (Other names include: (2-Aminopropyl)
11 Benzofuran).
12 (IIII) AB-PINACA
13 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
14 indazole-3-carboxamide).
15 (JJJJ) AB-FUBINACA
16 (N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-
17 1H-indazole-3-carboxamide).
18 (KKKK) ADB-PINACA
19 (N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
20 indazole-3-carboxamide).
21 (LLLL) Fluoro ADBICA (N-(1-Amino-3,3-
22 dimethyl-1-oxobutan-2-yl)-(fluoropentyl)-1H-indole-3-
23 carboxamide).
24 (MMMM) APDB (Other names include: -EMA,
25 -Desoxy-MDA, and (2-Aminopropyl)-2,3-
26 dihydrobenzofuran).
27 (NNNN) THJ-2201 (Other names include: AM2201 indazole
28 analog, Fluoropentyl-JWH-018 indazole, and
29 5-Fluoro-THJ-018).
30 (OOOO) AM 2201 benzimidazole analog (Other names
31 include: FUBIMINA, FTHJ, and (1-(5-fluoropentyl)-1H-
32 benzo[d]imidazol-2-yl)(naphthalene-1-yl)methanone).
33 (PPPP) MN-25 (Other names include: 7-methoxy-1-
34 [2-(4-morpholinyl)ethyl]-N-[1S, 2S, 4R)-1,3,3-
35 trimethylbicyclo[2.2.1]hept-2-yl]-1H-indole-3-carboxamide
36 and UR-12).
37 (QQQQ) FUB-PB-22 (Other names include:
38 Quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).
39 (RRRR) FUD-PB-22 (Other names include:
40 Naphthalen-1-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate).

- 1 (SSSS) 5-Fluoro-AB-PINACA (Other names include:
 2 AB-PINACA 5-fluoro analog and N-(1-amino-3-methyl-
 3 oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-
 4 carboxamide).
- 5 (TTTT) 4-MePPP (Other names include:
 6 4-methyl-alpha-pyrrolidinopropiophenone).
- 7 (UUUU) alpha-PBP (Other names include:
 8 Alpha-pyrrolidinobutiophenone).
- 9 (VVVV) AB-CHMINACA (Other names include:
 10 (N-[1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethy
 11 l)-1H-indazole-3-carboxamide).
- 12 (WWWW) Mexedrone
 13 (3-methoxy-2-(methylamino)-1-(p-tolyl)propan-1-one).
 14 (XXXX) MT-45,
 15 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine).
- 16 (YYYY) methyl 2-(1-(5-fluoropentyl)- 1H-indazole-3-
 17 carboxamido) -3,3-dimethylbutanoate [5F-ADB;
 18 5F-MDMB-PINACA].
- 19 (ZZZZ) methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-
 20 carboxamido)-3-methylbutanoate [5F-AMB].
- 21 (AAAAA) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)
 22 -1-(4-fluorobenzyl)- 1H-indazole-3-carboxamide
 23 [ADB-FUBINACA].
- 24 (BBBBB) N-(adamantan-1-yl)-1-(5-fluoropentyl)-
 25 1H-indazole-3- carboxamide [5F-APINACA, 5F-AKB48].
- 26 (CCCCC) methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
 27 carboxamido)-3,3-dimethylbutanoate [MDMB-CHMICA,
 28 MMB-CHMINACA].
- 29 (DDDDD) methyl 2-(1-(4-fluorobenzyl)-
 30 1H-indazole-3-carboxamido)- 3,3-dimethylbutanoate
 31 [MDMB-FUBINACA].
- 32 (EEEEE) N-(1-amino-3,3-dimethyl-1 -oxobutan-2-yl)-1-
 33 (cyclohexylmethyl)- 1H-indazole-3-carboxamide
 34 [MAB-CHMINACA and ADB-CHMINACA].
- 35 (FFFFF) Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-
 36 3-carboxamido)- 3-methylbutanoate [FUB-AMB,
 37 MMB-FUBINACA, AMB-FUBINACA].
- 38 (GGGGG) 3,4-dichloro-N-[(1dimethylamino)cyclohexylme
 39 thyl]benzamide) [AH7921].
- 40 (HHHHH) Naphthalen-1-yl 1-(5-fluoropentyl)-1

1 H-indole-3-carboxylate (trivial name: NM2201; CBL2201)
2 (IIIII) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1
3 H-indazole-3-carboxamide (trivial name:
4 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINACA;
5 4-CN-CUMYLBINACA; CUMYL-4CN-BINACA; SGT-78).
6 (JJJJJ) methyl 2-(1-(cyclohexylmethyl)-1
7 H-indole-3-carboxamido)-3-methylbutanoate (trivial names:
8 MMB-CHMICA, AMB-CHMICA).
9 (KKKKK) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1
10 H-pyrrolo[2,3-b]pyridine-3-carboxamide (trivial name:
11 5F-CUMYL-P7AICA).
12 (LLLLL) N-1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-p
13 entanone (N-ethylpentylone, ephylone).
14 (MMMMM) Synthetic cathinone, 1-(1,3-benzodioxol-5-yl)-2-
15 (ethylamino)-pentan-1-one (N-ethylpentylone, ephylone) and
16 its optical, positional, and geometric isomers, salts, and salts
17 of isomers.
18 (NNNNN) ethyl
19 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-
20 dimethylbutanoate (trivial name: 5F-EDMB-PINACA).
21 (OOOOO) methyl 2-(1-(5- fluoropentyl)-1H-indole-3-
22 carboxamido)-3,3-dimethylbutanoate (trivial name:
23 5F-MDMB-PICA).
24 (PPPPP) N- (adamantan- 1-yl)- 1-(4-fluorobenzyl)-
25 1H-indazole-3-carboxamide (trivial names: FUB-AKB48;
26 FUB-APINACA; AKB48 N- (4-FLUOROBENZYL)).
27 (QQQQQ) 1-(5- fluoropentyl)-N-(2-phenylpropan-2-yl)-
28 1H-indazole-3-carboxamide (trivial names:
29 5F-CUMYL-PINACA; SGT-25).
30 (RRRRR) (1-(4-fluorobenzyl)-1H-indol-3-
31 yl)(2,2,3,3-tetramethylcyclopropyl) methanone (trivial name:
32 FUB-144).
33 (SSSSS) 4F-MDMB-BINACA.
34 (TTTTT) N - e t h y l h e x e d r o n e
35 (2-(ethylamino)-1-phenylhexan-1-one).
36 (UUUUU) alpha-pyrrolidinohexanophenone
37 (1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one).
38 (VVVVV) alpha-pyrrolidinohexiophenone; trivial name:
39 a-PHP.
40 (WWWWW) 4'-methyl-alpha-pyrrolidinohexiophenone

- 1 (1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one.
 2 (XXXXXX) 4-methyl-alphaethylaminopentiophenone
 3 (2-(ethylamino)-1-(4-methylphenyl)pentan-1-one; trivial
 4 name: 4-MEAP.
 5 (YYYYYY) 4'-methyl-alphapyrrolidinohexanophenone; trivial
 6 name: MPHP.
 7 (ZZZZZ) alphapyrrolidinoheptaphenone
 8 (1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one; trivial name: PV8.
 9 (AAAAAA) 4'-chloro-alphapyrrolidinovalerophenone (1-(4-
 10 chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one.
 11 (BBBBBB) 4'-chloro-alphapyrrolidinopentiophenone; trivial
 12 name: 4-chloro-a-PVP.
 13 **(CCCCCC) 4,4'-dimethylaminorex (common name:**
 14 **4,4'-DMAR).**
 15 (2) Any compound structurally derived from
 16 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by
 17 substitution at the nitrogen atom of the indole ring by alkyl,
 18 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
 19 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 20 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 21 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 22 or not further substituted in the indole ring to any extent and
 23 whether or not substituted in the naphthyl ring to any extent.
 24 (3) Any compound structurally derived from 3-(1-naphthoyl)
 25 pyrrole by substitution at the nitrogen atom of the pyrrole ring by
 26 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
 27 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,
 28 2-(4-morpholinyl)ethyl, or 1-(N-methyl-2-pyrrolidinyl)methyl,
 29 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
 30 group, whether or not further substituted in the pyrrole ring to any
 31 extent and whether or not substituted in the naphthyl ring to any
 32 extent.
 33 (4) Any compound structurally derived from
 34 1-(1-naphthylmethyl)indene by substitution at the 3-position of
 35 the indene ring by alkyl, haloalkyl, cyanoalkyl, alkenyl,
 36 c y c l o a l k y l m e t h y l , c y c l o a l k y l e t h y l ,
 37 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, or
 38 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
 39 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
 40 or not further substituted in the indene ring to any extent and

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

1 (9) Any compound structurally derived from 3-tetramethyl
2 cyclopropanoylindole with substitution at the nitrogen atom of the
3 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
4 c y c l o a l k y l m e t h y l , c y c l o a l k y l e t h y l ,
5 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl) ethyl,
6 1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
7 morpholinyl)methyl, or tetrahydropyranylmethyl group, whether
8 or not further substituted in the indole ring to any extent and
9 whether or not substituted in the tetramethylcyclopropyl ring to
10 any extent.

11 (10) Any compound containing a N-(1-adamantyl)-
12 1H-indazole-3-carboxamide structure with substitution at the
13 nitrogen atom of the indazole ring by an alkyl, haloalkyl,
14 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
15 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
16 1 - (N - m e t h y l - 2 - p y r r o l i d i n y l) m e t h y l ,
17 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
18 group, whether or not further substituted at the nitrogen atom of
19 the carboxamide to any extent, whether or not further substituted
20 in the indazole ring to any extent, and whether or not further
21 substituted on the adamantyl ring system to any extent. An
22 example of this structural class includes AKB48.

23 (11) Any compound containing a N-(1-adamantyl)-
24 1H-indole-3-carboxamide structure with substitution at the
25 nitrogen atom of the indole ring by an alkyl, haloalkyl,
26 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
27 1-(N-methyl-2- piperidinyl)methyl, or 2-(4-morpholinyl)ethyl,
28 1 - (N - m e t h y l - 2 - p y r r o l i d i n y l) m e t h y l ,
29 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl
30 group, whether or not further substituted at the nitrogen atom of
31 the carboxamide to any extent, whether or not further substituted
32 in the indole ring to any extent, and whether or not further
33 substituted on the adamantyl ring system to any extent. An
34 example of this structural class includes STS-135.

35 (12) Any compound containing a 3-(1-adamantoyl)indole
36 structure with substitution at the nitrogen atom of the indole ring
37 by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
38 cycloalkylethyl, 1-(N-methyl-2- piperidinyl)methyl, or
39 2-(4-morpholinyl)ethyl, 1-(N-methyl-2- pyrrolidinyl)methyl,
40 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl

1 group, whether or not further substituted on the adamantyl ring
 2 system to any extent. An example of this structural class includes
 3 AM-1248.

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

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

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

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

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

17 **"Brorphine".**

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

19 **"2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-**
 20 **benzimidazole (N-pyrrolidino etonitazene; etonitazepyne)".**

21 Page 3, delete line 1.

22 Page 6, strike line 1.

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

25 Renumber all SECTIONS consecutively.

(Reference is to SB 209 as introduced.)