HB5268

LRB10119113RLC68576b

101ST GENERAL ASSEMBLY

State of Illinois

2019 and 2020

HB5268

by Rep. John Connor

SYNOPSIS AS INTRODUCED:

720 ILCS 570/204 from Ch. 56 1/2, par. 1204

Amends the Illinois Controlled Substances Act. Includes various fentanyl analogs in the list of Schedule I controlled substances.

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CORRECTIONAL
BUDGET AND
IMPACT NOTE ACT
MAY APPLY

A BILL FOR

AN ACT concerning criminal law.

Be it enacted by the People of the State of Illinois, represented in the General Assembly:

Section 5. The Illinois Controlled Substances Act is amended by changing Section 204 as follows:

(720 ILCS 570/204) (from Ch. 56 1/2, par. 1204)

Sec. 204. (a) The controlled substances listed in this Section are included in Schedule I.

(b) Unless specifically excepted or unless listed in another schedule, any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers and salts is possible within the specific chemical designation:

(1) Acetylmethadol;

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

(2) Allylprodine;

(3) Alphacetylmethadol, except levo-alphacetylmethadol (also known as levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM);

(4) Alphameprodine;

(5) Alphamethadol;

(6) Alpha-methylfentanyl (N-(1-alpha-methyl-beta-phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine;

(6.1) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

(7) 1-methyl-4-phenyl-4-propionoxypiperidine (MPPP);

(7.1) PEPAP (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);

(8) Benzethidine;

(9) Betacetylmethadol;

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

(10) Betameprodine;

(11) Betamethadol;

(12) Betaprodine;

(13) Clonitazene;

(14) Dextromoramide;

(15) Diampromide;

(16) Diethylthiambutene;

(17) Difenoxin;

(18) Dimenoxadol;

(19) Dimepheptanol;

(20) Dimethylthiambutene;

(21) Dioxaphetylbutyrate;

(22) Dipipanone;

(23) Ethylmethylthiambutene;

(24) Etonitazene;

(25) Etoxeridine;

(26) Furethidine;

(27) Hydroxpethidine;

(28) Ketobemidone;

(29) Levomoramide;

(30) Levophenacylmorphan;

(31) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

(31.1) 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

(32) Morpheridine;

(33) Noracymethadol;

(34) Norlevorphanol;

(35) Normethadone;

(36) Norpipanone;

(36.1) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl]propanamide);

(37) Phenadoxone;

(38) Phenampromide;

(39) Phenomorphan;

(40) Phenoperidine;

(41) Piritramide;

(42) Proheptazine;

(43) Properidine;

(44) Propiram;

(45) Racemoramide;

(45.1) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);

(46) Tilidine;

(47) Trimeperidine;

(48) Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide);

(49) Furanyl fentanyl (FU-F);

(50) Butyryl fentanyl;

(51) Valeryl fentanyl;

(52) Acetyl fentanyl;

(53) Beta-hydroxy-thiofentanyl;

(54) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (U-47700);

(55) 4-chloro-N-[1-[2-(4-nitrophenyl)ethyl]-2-piperidinylidene]-

benzenesulfonamide (W-18);

(56) 4-chloro-N-[1-(2-phenylethyl) -2-piperidinylidene]-benzenesulfonamide (W-15);

(57) acrylfentanyl (acryloylfentanyl)~~.~~;

(b-5) Fentanyl analogs, including any compound, except compounds scheduled elsewhere in this Act, structurally derived from fentanyl by replacement of the phenyl portion of the phenethyl group by any monocycle whether or not further substituted in or on the monocycle; by substitution in or on the phenethyl group with alkyl, alkenyl, alkoxy, hydroxy, halo, haloalkyl, amino or nitro groups; by substitution in or on the piperidine ring with alkyl, alkenyl, alkoxy, ester, ether, hydroxy, halo, haloalkyl, amino or nitro groups; by replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; by replacement of the N-propionyl group by another acyl group; or by any combination of these modifications. Substances specified under this subsection include all of the following:

(1) Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);

(2) 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl) piperidin-4-yl]-butyramide);

(3) Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]

-propanamide);

(4) Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);

(5) 3-methylbutyryl fentanyl(N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-phenylbutyramide);

(6) 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);

(7) Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);

(8) Tetrahydrofuran fentanyl(N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide);

(9) Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).

(c) Unless specifically excepted or unless listed in another schedule, any of the following opium derivatives, 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:

(1) Acetorphine;

(2) Acetyldihydrocodeine;

(3) Benzylmorphine;

(4) Codeine methylbromide;

(5) Codeine-N-Oxide;

(6) Cyprenorphine;

(7) Desomorphine;

(8) Diacetyldihydromorphine (Dihydroheroin);

(9) Dihydromorphine;

(10) Drotebanol;

(11) Etorphine (except hydrochloride salt);

(12) Heroin;

(13) Hydromorphinol;

(14) Methyldesorphine;

(15) Methyldihydromorphine;

(16) Morphine methylbromide;

(17) Morphine methylsulfonate;

(18) Morphine-N-Oxide;

(19) Myrophine;

(20) Nicocodeine;

(21) Nicomorphine;

(22) Normorphine;

(23) Pholcodine;

(24) Thebacon.

(d) Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following hallucinogenic substances, or which contains any of 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 (for the purposes of this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

(1) 3,4-methylenedioxyamphetamine (alpha-methyl,3,4-methylenedioxyphenethylamine, methylenedioxyamphetamine, MDA);

(1.1) Alpha-ethyltryptamine (some trade or other names: etryptamine; MONASE; alpha-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)indole; a-ET; and AET);

(2) 3,4-methylenedioxymethamphetamine (MDMA);

(2.1) 3,4-methylenedioxy-N-ethylamphetamine (also known as: N-ethyl-alpha-methyl-3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE, and MDEA);

(2.2) N-Benzylpiperazine (BZP);

(2.2-1) Trifluoromethylphenylpiperazine (TFMPP);

(3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);

(4) 3,4,5-trimethoxyamphetamine (TMA);

(5) (Blank);

(6) Diethyltryptamine (DET);

(7) Dimethyltryptamine (DMT);

(7.1) 5-Methoxy-diallyltryptamine;

(8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);

(9) Ibogaine (some trade and other names:

7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b] indole; Tabernanthe iboga);

(10) Lysergic acid diethylamide;

(10.1) Salvinorin A;

(10.5) Salvia divinorum (meaning all parts of the plant presently classified botanically as Salvia divinorum, whether growing or not, the seeds thereof, any extract from any part of that plant, and every compound, manufacture, salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation, derivative, mixture, or preparation of that plant, its seeds or extracts);

(11) 3,4,5-trimethoxyphenethylamine (Mescaline);

(12) Peyote (meaning all parts of the plant presently classified botanically as Lophophora williamsii Lemaire, whether growing or not, the seeds thereof, any extract from any part of that plant, and every compound, manufacture, salts, derivative, mixture, or preparation of that plant, its seeds or extracts);

(13) N-ethyl-3-piperidyl benzilate (JB 318);

(14) N-methyl-3-piperidyl benzilate;

(14.1) N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine and N-hydroxy MDA);

(15) Parahexyl; some trade or other names: 3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo (b,d) pyran; Synhexyl;

(16) Psilocybin;

(17) Psilocyn;

(18) Alpha-methyltryptamine (AMT);

(19) 2,5-dimethoxyamphetamine (2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA);

(20) 4-bromo-2,5-dimethoxyamphetamine (4-bromo-2,5-dimethoxy-alpha-methylphenethylamine; 4-bromo-2,5-DMA);

(20.1) 4-Bromo-2,5 dimethoxyphenethylamine. Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl DOB, 2CB, Nexus;

(21) 4-methoxyamphetamine (4-methoxy-alpha-methylphenethylamine; paramethoxyamphetamine; PMA);

(22) (Blank);

(23) Ethylamine analog of phencyclidine. Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE;

(24) Pyrrolidine analog of phencyclidine. Some trade or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy,

PHP;

(25) 5-methoxy-3,4-methylenedioxy-amphetamine;

(26) 2,5-dimethoxy-4-ethylamphetamine (another name: DOET);

(27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine (another name: TCPy);

(28) (Blank);

(29) Thiophene analog of phencyclidine (some trade or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 2-thienyl analog of phencyclidine; TPCP; TCP);

(29.1) Benzothiophene analog of phencyclidine. Some trade or other names: BTCP or benocyclidine;

(29.2) 3-Methoxyphencyclidine (3-MeO-PCP);

(30) Bufotenine (some trade or other names: 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; 5-hydroxy-N,N-dimethyltryptamine; N,N-dimethylserotonin; mappine);

(31) (Blank);

(32) (Blank);

(33) (Blank);

(34) (Blank);

(34.5) (Blank);

(35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol

Some trade or other names: HU-210;

(35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol, its isomers, salts, and salts of isomers; Some trade or other names: HU-210, Dexanabinol;

(36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol

Some trade or other names: HU-211;

(37) (Blank);

(38) (Blank);

(39) (Blank);

(40) (Blank);

(41) (Blank);

(42) 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, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl whether or not further substituted in the indole ring to any extent, whether or not substituted in the naphthyl ring to any extent. Examples of this structural class include, but are not limited to, JWH-018, AM-2201, JWH-175, JWH-184, and JWH-185;

(43) Any compound structurally derived from 3-(1-naphthoyl)pyrrole by substitution at the nitrogen atom of the pyrrole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted in the pyrrole ring to any extent, whether or not substituted in the naphthyl ring to any extent. Examples of this structural class include, but are not limited to, JWH-030, JWH-145, JWH-146, JWH-307, and JWH-368;

(44) Any compound structurally derived from 1-(1-naphthylmethyl)indene by substitution at the 3-position of the indene ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl whether or not further substituted in the indene ring to any extent, whether or not substituted in the naphthyl ring to any extent. Examples of this structural class include, but are not limited to, JWH-176;

(45) Any compound structurally derived from 3-phenylacetylindole by substitution at the nitrogen atom of the indole ring with alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted

in the indole ring to any extent, whether or not substituted in the phenyl ring to any extent. Examples of this structural class include, but are not limited to, JWH-167, JWH-250, JWH-251, and RCS-8;

(46) Any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not substituted in the cyclohexyl ring to any extent. Examples of this structural class include, but are not limited to, CP 47, 497 and its C8 homologue (cannabicyclohexanol);

(46.1) Any compound structurally derived from 3-(benzoyl) indole with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl 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. Examples of this structural class include, but are not limited to, AM-630, AM-2233, AM-694, Pravadoline (WIN 48,098), and RCS-4;

(47) (Blank);

(48) (Blank);

(49) (Blank);

(50) (Blank);

(51) (Blank);

(52) (Blank);

(53) 2,5-Dimethoxy-4-(n)-propylthio-phenethylamine. Some trade or other names: 2C-T-7;

(53.1) 4-ethyl-2,5-dimethoxyphenethylamine. Some trade or other names: 2C-E;

(53.2) 2,5-dimethoxy-4-methylphenethylamine. Some trade or other names: 2C-D;

(53.3) 4-chloro-2,5-dimethoxyphenethylamine. Some trade or other names: 2C-C;

(53.4) 4-iodo-2,5-dimethoxyphenethylamine. Some trade or other names: 2C-I;

(53.5) 4-ethylthio-2,5-dimethoxyphenethylamine. Some trade or other names: 2C-T-2;

(53.6) 2,5-dimethoxy-4-isopropylthio-phenethylamine. Some trade or other names: 2C-T-4;

(53.7) 2,5-dimethoxyphenethylamine. Some trade or other names: 2C-H;

(53.8) 2,5-dimethoxy-4-nitrophenethylamine. Some trade or other names: 2C-N;

(53.9) 2,5-dimethoxy-4-(n)-propylphenethylamine. Some trade or other names: 2C-P;

(53.10) 2,5-dimethoxy-3,4-dimethylphenethylamine. Some trade or other names: 2C-G;

(53.11) The N-(2-methoxybenzyl) derivative of any 2C phenethylamine referred to in subparagraphs (20.1), (53), (53.1), (53.2), (53.3), (53.4), (53.5), (53.6), (53.7), (53.8), (53.9), and (53.10) including, but not limited to, 25I-NBOMe and 25C-NBOMe;

(54) 5-Methoxy-N,N-diisopropyltryptamine;

(55) (Blank);

(56) (Blank);

(57) (Blank);

(58) (Blank);

(59) 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any extent: including, but not limited to, XLR11, UR144, FUB-144;

(60) 3-adamantoylindole with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any extent: including, but not limited to, AB-001;

(61) N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any extent: including, but not limited to, APICA/2NE-1, STS-135;

(62) N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the adamantyl ring to any extent: including, but not limited to, AKB48, 5F-AKB48;

(63) 1H-indole-3-carboxylic acid 8-quinolinyl ester with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent, whether or not substituted on the quinoline ring to any extent: including,

but not limited to, PB22, 5F-PB22, FUB-PB-22;

(64) 3-(1-naphthoyl)indazole with substitution at the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the naphthyl ring to any extent: including, but not limited to, THJ-018, THJ-2201;

(65) 2-(1-naphthoyl)benzimidazole with substitution at the nitrogen atom of the benzimidazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the benzimidazole ring to any extent, whether or not substituted on the naphthyl ring to any extent: including, but not limited to, FUBIMINA;

(66) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, AB-PINACA, AB-FUBINACA, AB-CHMINACA;

(67) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, ADB-PINACA, ADB-FUBINACA;

(68) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent: including, but not limited to, ADBICA, 5F-ADBICA;

(69) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indole ring to any extent: including, but not limited to, ABICA, 5F-ABICA;

(70) Methyl 2-(1H-indazole-3-carboxamido)-3-methylbutanoate with substitution on the nitrogen atom of

the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, AMB, 5F-AMB;

(71) Methyl 2-(1H-indazole-3-carboxamido)-3,3-dimethylbutanoate with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, 5-fluoro-MDMB-PINACA, MDMB-FUBINACA;

(72) Methyl 2-(1H-indole-3-carboxamido)-3-methylbutanoate with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, MMB018, MMB2201, and AMB-CHMICA;

(73) Methyl 2-(1H-indole-3-carboxamido)-3,3-dimethylbutanoate with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl

halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, MDMB-CHMICA;

(74) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-indazole-3-carboxamide with substitution on the nitrogen atom of the indazole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, APP-CHMINACA, 5-fluoro-APP-PINACA;

(75) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1H-indole-3-carboxamide with substitution on the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, aryl halide, alkyl aryl halide, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl, whether or not further substituted on the indazole ring to any extent: including, but not limited to, APP-PICA and 5-fluoro-APP-PICA;

(76) 4-Acetoxy-N,N-dimethyltryptamine: trade name 4-AcO-DMT;

(77) 5-Methoxy-N-methyl-N-isopropyltryptamine: trade name 5-MeO-MIPT;

(78) 4-hydroxy Diethyltryptamine (4-HO-DET);

(79) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET);

(80) 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT);

(81) 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT);

(82) Fluorophenylpiperazine;

(83) Methoxetamine;

(84) 1-(Ethylamino)-2-phenylpropan-2-one (iso-ethcathinone).

(e) Unless specifically excepted 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:

(1) mecloqualone;

(2) methaqualone; and

(3) gamma hydroxybutyric acid.

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

(1) Fenethylline;

(2) N-ethylamphetamine;

(3) Aminorex (some other names: 2-amino-5-phenyl-2-oxazoline; aminoxaphen;

4-5-dihydro-5-phenyl-2-oxazolamine) and its salts, optical isomers, and salts of optical isomers;

(4) Methcathinone (some other names: 2-methylamino-1-phenylpropan-1-one; Ephedrone; 2-(methylamino)-propiophenone; alpha-(methylamino)propiophenone; N-methylcathinone; methycathinone; Monomethylpropion; UR 1431) and its salts, optical isomers, and salts of optical isomers;

(5) Cathinone (some trade or other names: 2-aminopropiophenone; alpha-aminopropiophenone; 2-amino-1-phenyl-propanone; norephedrone);

(6) N,N-dimethylamphetamine (also known as: N,N-alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine);

(7) (+ or -) cis-4-methylaminorex ((+ or -) cis-4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);

(8) 3,4-Methylenedioxypyrovalerone (MDPV);

(9) Halogenated amphetamines and methamphetamines - any compound derived from either amphetamine or methamphetamine through the substitution of a halogen on the phenyl ring, including, but not limited to, 2-fluoroamphetamine, 3-fluoroamphetamine and 4-fluoroamphetamine;

(10) Aminopropylbenzofuran (APB): including 4-(2-Aminopropyl) benzofuran, 5-(2-Aminopropyl)benzofuran, 6-(2-Aminopropyl)

benzofuran, and 7-(2-Aminopropyl) benzofuran;

(11) Aminopropyldihydrobenzofuran (APDB): including 4-(2-Aminopropyl)-2,3- dihydrobenzofuran, 5-(2-Aminopropyl)-2, 3-dihydrobenzofuran, 6-(2-Aminopropyl)-2,3-dihydrobenzofuran, and 7-(2-Aminopropyl)-2,3-dihydrobenzofuran;

(12) Methylaminopropylbenzofuran (MAPB): including 4-(2-methylaminopropyl) benzofuran, 5-(2-methylaminopropyl)benzofuran, 6-(2-methylaminopropyl)benzofuran and 7-(2-methylaminopropyl)benzofuran.

(g) Temporary listing of substances subject to emergency scheduling. Any material, compound, mixture, or preparation that contains any quantity of the following substances:

(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, isomers, salts, and salts of isomers;

(2) N-[1(2-thienyl) methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical isomers, salts, and salts of isomers.

(h) Synthetic cathinones. Unless specifically excepted, any chemical compound which is not approved by the United States Food and Drug Administration or, if approved, is not dispensed or possessed in accordance with State or federal law, not including bupropion, 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 in one or more of the following ways:

(1) 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. Examples of this class include, but are not limited to, 3,4-Methylenedioxycathinone (bk-MDA);

(2) by substitution at the 3-position with an acyclic alkyl substituent. Examples of this class include, but are not limited to, 2-methylamino-1-phenylbutan-1-one (buphedrone); or

(3) by substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or methoxybenzyl groups, or by inclusion of the 2-amino nitrogen atom in a cyclic structure. Examples of this class include, but are not limited to, Dimethylcathinone, Ethcathinone, and a-Pyrrolidinopropiophenone (a-PPP); or

Any other synthetic cathinone which is not approved by the United States Food and Drug Administration or, if approved, is not dispensed or possessed in accordance with State or federal law.

(i) Synthetic cannabinoids or piperazines. Any synthetic cannabinoid or piperazine which is not approved by the United States Food and Drug Administration or, if approved, which is

not dispensed or possessed in accordance with State and federal law.

(Source: P.A. 99-371, eff. 1-1-16; 100-201, eff. 8-18-17; 100-368, eff. 1-1-18; 100-789, eff. 1-1-19; 100-863, eff. 8-14-18.)