HB2526

WEST VIRGINIA LEGISLATURE

2017 REGULAR SESSION

ENROLLED

Committee Substitute

for

House Bill 2526

BY DELEGATES ELLINGTON, SUMMERS, SOBONYA AND

ROHRBACH

[Passed April 8, 2017; in effect ninety days from passage.]

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AN ACT to amend and reenact §60A-2-201, §60A-2-204, §60A-2-206, §60A-2-210 and §60A-2-212 of the Code of West Virginia, 1931, as amended, all relating to classifying additional drugs to Schedules I, II, IV and V of controlled substances; and adding a provision relating to the scheduling of a cannabidiol in a product approved by the Food and Drug Administration.

Be it enacted by the Legislature of West Virginia:

That §60A-2-201, §60A-2-204, §60A-2-206, §60A-2-210 and §60A-2-212 of the Code of West Virginia, 1931, as amended, be amended and reenacted, all to read as follows:

ARTICLE 2. STANDARDS AND SCHEDULES.

§60A-2-201. Authority of state Board of Pharmacy; recommendations to Legislature.

(a) The state Board of Pharmacy shall administer the provisions of this chapter. It shall also, on the first day of each regular legislative session, recommend to the Legislature which substances should be added to or deleted from the schedules of controlled substances contained in this article or reschedule therein. The state Board of Pharmacy shall also have the authority between regular legislative sessions, on an emergency basis, to add to or delete from the schedules of controlled substances contained in this article or reschedule such substances based upon the recommendations and approval of the federal food, drug and cosmetic agency, and shall report such actions on the first day of the regular legislative session immediately following said actions.

In making any such recommendation regarding a substance, the state Board of Pharmacy shall consider the following factors:

(1) The actual or relative potential for abuse;

(2) The scientific evidence of its pharmacological effect, if known;

(3) The state of current scientific knowledge regarding the substance;

(4) The history and current pattern of abuse;

(5) The scope, duration and significance of abuse;

(6) The potential of the substance to produce psychic or physiological dependence liability; and

(7) Whether the substance is an immediate precursor of a substance already controlled under this article.

(b) After considering the factors enumerated in subsection (a), the state Board of Pharmacy shall make findings with respect to the substance under consideration. If it finds that any substance not already controlled under any schedule has a potential for abuse, it shall recommend to the Legislature that the substance be added to the appropriate schedule. If it finds that any substance already controlled under any schedule should be rescheduled or deleted, it shall so recommend to the Legislature.

(c) If the state Board of Pharmacy designates a substance as an immediate precursor, substances which are precursors of the controlled precursor shall not be subject to control solely because they are precursors of the controlled precursor.

(d) If any substance is designated, rescheduled or deleted as a controlled substance under federal laws and notice thereof is given to the state Board of Pharmacy, the board shall recommend similar control of such substance to the Legislature, specifically stating that such recommendation is based on federal action and the reasons why the federal government deemed such action necessary and proper.

(e) The authority vested in the board by subsection (a) of this section shall not extend to distilled spirits, wine, malt beverages or tobacco as those terms are defined or used in other chapters of this code nor to any nonnarcotic substance if such substance may under the "Federal Food, Drug and Cosmetic Act" and the law of this state lawfully be sold over the counter without a prescription.

(f) Notwithstanding any provision of this chapter to the contrary, the sale, wholesale, distribution or prescribing of a cannabidiol in a product approved by the Food and Drug Administration is permitted and shall be placed on the schedule as provided for by the Drug Enforcement Administration.

60A-2-204. Schedule I.

(a) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.

(b) Opiates. 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 (for purposes of subdivision (34) of this subsection only, the term isomer includes the optical and geometric isomers):

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

(2) Acetylmethadol;

(3) Allylprodine;

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

(5) Alphameprodine;

(6) Alphamethadol;

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

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

(9) Benzethidine;

(10) Betacetylmethadol;

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

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

(13) Betameprodine;

(14) Betamethadol;

(15) Betaprodine;

(16) Clonitazene;

(17) Dextromoramide;

(18) Diampromide;

(19) Diethylthiambutene;

(20) Difenoxin;

(21) Dimenoxadol;

(22) Dimepheptanol;

(23) Dimethylthiambutene;

(24) Dioxaphetyl butyrate;

(25) Dipipanone;

(26) Ethylmethylthiambutene;

(27) Etonitazene;

(28) Etoxeridine;

(29) Furethidine;

(30) Hydroxypethidine;

(31) Ketobemidone;

(32) Levomoramide;

(33) Levophenacylmorphan;

(34) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4- piperidyl]-N-phenylpropanamide);

(35) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4- piperidinyl]—phenylpropanamide);

(36) Morpheridine;

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

(38) Noracymethadol;

(39) Norlevorphanol;

(40) Normethadone;

(41) Norpipanone;

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

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

(44) Phenadoxone;

(45) Phenampromide;

(46) Phenomorphan;

(47) Phenoperidine;

(48) Piritramide;

(49) Proheptazine;

(50) Properidine;

(51) Propiram;

(52) Racemoramide;

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

(54) Tilidine;

(55) Trimeperidine.

(c) *Opium derivatives*. — Unless specifically excepted or unless listed in another schedule, any of the following opium immediate 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) Dihydromorphine;

(9) Drotebanol;

(10) Etorphine (except HCl Salt);

(11) Heroin;

(12) Hydromorphinol;

(13) Methyldesorphine;

(14) Methyldihydromorphine;

(15) Morphine methylbromide;

(16) Morphine methylsulfonate;

(17) Morphine-N-Oxide;

(18) Myrophine;

(19) Nicocodeine;

(20) Nicomorphine;

(21) Normorphine;

(22) Pholcodine;

(23) Thebacon.

(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 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 purposes of this subsection only, the term "isomer" includes the optical, position and geometric isomers):

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

(2) 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;

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

(4)(A) N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the acronym 25B-NBOMe.

(B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe).

(C) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)

(5) 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA;

(6) 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

(7) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

(8) 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-methylphenethylamine; paramethoxyamphetamine; PMA;

(9) 5-methoxy-3, 4-methylenedioxy-amphetamine;

(10) 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

(11) 3,4-methylenedioxy amphetamine;

(12) 3,4-methylenedioxymethamphetamine (MDMA);

(13) 3,4-methylenedioxy-N-ethylamphetamine (also known as – ethyl-alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

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

(15) 3,4,5-trimethoxy amphetamine;

(16) 5-methoxy-N, N-dimethyltryptamine (5-MeO-DMT);

(17) Alpha-methyltryptamine (other name: AMT);

(18) Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;

(19) Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;

(20) Dimethyltryptamine; some trade or other names: DMT;

(21) 5-Methoxy-N, N-diisopropyltryptamine (5-MeO-DIPT);

(22) 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;

(23) Lysergic acid diethylamide;

(24) Marijuana;

(25) Mescaline;

(26) 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; Synhexyl;

(27) 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 such plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of such plant, its seeds or extracts;

(28) N-ethyl-3-piperidyl benzilate;

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

(30) Psilocybin;

(31) Psilocyn;

(32) Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives and their isomers with similar chemical structure and pharmacological activity such as the following:

delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;

delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;

delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;

(Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions covered).

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

(34) Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;

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

(36) 1[1-(2-thienyl)cyclohexyl]pyrroldine; some other names: TCPy.

(37) 4-methylmethcathinone (Mephedrone);

(38) 3,4-methylenedioxypyrovalerone (MDPV);

(39) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

(40) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

(41) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

(42) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

(43) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

(44) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

(45) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

(46) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N);

(47) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);

(48) 3,4-Methylenedioxy-N-methylcathinone (Methylone);

(49) (2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, itsoptical isomers, salts and salts of isomers

(50) 5-methoxy-N, N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);

(51) Alpha-methyltryptamine (other name: AMT);

(52) 5-methoxy-N, N-diisopropyltryptamine (other name: 5-MeO-DIPT);

(53) Synthetic Cannabinoids as follows:

(A) 2-[(1R,3S)-3-hydroxycyclohexyl]-5- (2-methyloctan-2-yl) phenol) {also known as CP 47,497 and homologues};

(B) rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl) phenol {also known as CP 47,497-C8 homolog};

(C) [(6a*R*)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7,10,10a-tetrahydrobenzo[c]chromen-1-ol)] {also known as HU-210};

(D) (dexanabinol);

(6a*S*,10a*S*)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzol[c]chromen-1-ol) {also known as HU-211};

(E) 1-Pentyl-3-(1-naphthoyl) indole {also known as JWH-018};

(F) 1-Butyl-3-(1-naphthoyl) indole {also known as JWH-073};

(G) (2-methyl-1-propyl-1H-indol-3-yl)-1-napthalenyl-methanone {also known as JWH-015};

(H) (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};

(I) [1-[2-(4-morpholinyl) ethyl] -1H-indol-3-yl]-1-naphthalenyl-methanone {also known as JWH-200};

(J) 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};

(K) 2-((1S,2S,5S)-5-hydroxy-2- (3-hydroxtpropyl)cyclohexyl) -5-(2-methyloctan-2-yl)phenol {also known as CP 55,940};

(L) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-122};

(M) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-398;

(N) (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};

(O) 1-(1-(2-cyclohexylethyl) -1H-indol-3-yl) -2-(2-methoxyphenyl) ethanone {also known as RCS-8};

(P) 1-pentyl-3-[1-(4-methoxynaphthoyl) indole (JWH-081);

(Q) 1-(5-fluoropentyl)-3-(1-naphthoyl) indole (AM2201); and

(R) 1-(5-fluoropentyl)-3-(2-iodobenzoyl) indole (AM694).

(54) Synthetic cannabinoids or any material, compound, mixture or preparation which contains any quantity of the following substances, including their analogues, congeners, homologues, isomers, salts and salts of analogues, congeners, homologues and isomers, as follows:

(A) CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-YL) phenol);

(B) HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10, 10A-tetrahydrobenzo[C] chromen-1-OL)];

(C) HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-YL)-6A,7,10,10 atetrahydrobenzo [C] chromen-1-OL);

(D) JWH-018, 1-pentyl-3-(1-naphthoyl) indole;

(E) JWH-019, 1-hexyl-3-(1-naphthoyl) indole;

(F) JWH-073, 1-butyl-3-(1-naphthoyl) indole;

(G) JWH-200, (1-(2-morpholin-4-ylethyl) indol-3-yl)- Naphthalen-1-ylmethanone;

(H) JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl) indole.

(55) Synthetic cannabinoids including any material, compound, mixture or preparation that is not listed as a controlled substance in Schedule I through V, is not a federal Food and Drug Administration approved drug or used within legitimate and approved medical research and which contains any quantity of the following substances, their salts, isomers, whether optical positional or geometric, analogues, homologues and salts of isomers, analogues and homologues, unless specifically exempted, whenever the existence of these salts, isomers, analogues, homologues and salts of isomers, analogues and homologues if possible within the specific chemical designation:

(A) Tetrahydrocannabinols meaning tetrahydrocannabinols which are naturally contained in a plant of the genus cannabis as well as synthetic equivalents of the substances contained in the plant or in the resinous extractives of cannabis or synthetic substances, derivatives and their isomers with analogous chemical structure and or pharmacological activity such as the following:

(i) DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.

(ii) DELTA-6 CIS OR trans tetrahydrocannabinol and their Optical isomers.

(iii) DELTA-3,4 CIS OR their trans tetrahydrocannabinol and their optical isomers.

(B) Naphthoyl indoles or any compound containing a 3-(-1- Napthoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include the following:

(i) JWH 015;

(ii) JWH 018;

(iii) JWH 019;

(iv) JWH 073;

(v) JWH 081;

(vi) JWH 122;

(vii) JWH 200;

(viii) JWH 210;

(ix) JWH 398;

(x) AM 2201;

(xi) WIN 55,212.

(56) Synthetic Phenethylamines (including their optical, positional, and geometric isomers, salts and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers):

(A) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-NBOMe);

(B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-NBOMe);

(C) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-NBOMe);

(57) Synthetic Opioids (icluding their isomers, esters, ethers, salts and salts of isomers, esters and ethers):

(A) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);

(B) furanyl fentanyl;

(C) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-47700);

(D) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);

(E) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethylpiperidin-4-yl]-N-phenylpropionamide, also known as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-hydroxythiofentanyl).

(58) Opioid Receptor Agonist (including its isomers, esters, ethers, salts, and salts of isomers, esters and ethers):

(A) AH-7921 (3,4-dichloro-N- (1dimethylamino)cyclohexylmethyl]benzamide).

(59) Naphylmethylindoles or any compound containing a 1hindol-3-yl-(1-naphthyl) methane structure with a substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 175 and JWH 184.

(60) Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure with substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the pyrrole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 147 and JWH 307.

(61) Naphthylmethylindenes or any compound containing a Naphthylideneindene structure with substitution at the 3- Position of the indene ring whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 176.

(62) Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include the following:

(A) RCS-8, SR-18 OR BTM-8;

(B) JWH 250;

(C) JWH 203;

(D) JWH 251;

(E) JWH 302.

(63) Cyclohexylphenols or any compound containing a 2-(3- hydroxycyclohexyl) phenol structure with a substitution at the 5-position of the phenolic ring whether or not substituted in the cyclohexyl ring to any extent. This shall include the following:

(A) CP 47,497 and its homologues and analogs;

(B) Cannabicyclohexanol;

(C) CP 55,940.

(64) Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with substitution at the nitrogren atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include the following:

(A) AM 694;

(B) Pravadoline WIN 48,098;

(C) RCS 4;

(D) AM 679.

(65) [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-napthalenymethanone. This shall include WIN 55,212-2.

(66) Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051 and JWH 133.

(67) Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the adamantoyl ring system to any extent. This shall include AM1248.

(68) Tetramethylcyclopropylindoles or any compound containing A 3-tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring

whether or not further substituted in the indole ring to any extent and whether or not substituted in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.

(69) N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.

(70) Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and V, not federal Food and Drug Administration approved drug or used within legitimate, approved medical research. Since nomenclature of these substances is not internationally standardized, any immediate precursor or immediate derivative of these substances shall be covered.

(71) Tryptamines:

(A) 5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)

(B) 4-hydroxy-N, N-diisopropyltryptamine (4-HO-DiPT)

(C) 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)

(D) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)

(E) 4-acetoxy-N, N-diisopropyltryptamine (4-AcO-DiPT)

(F) 5-methoxy-α-methyltryptamine (5-MeO-AMT)

(G) 4-methoxy-N, N-Dimethyltryptamine (4-MeO-DMT)

(H) 4-hydroxy Diethyltryptamine (4-HO-DET)

(I) 5- methoxy- N, N- diallyltryptamine (5-MeO-DALT)

(J) 4-acetoxy-N, N-Dimethyltryptamine (4-AcO DMT)

(K) 4-hydroxy Diethyltryptamine (4-HO-DET)

(72) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (AB-CHMINACA);

(73) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (AB-PINACA);

(74) [1-(5-fluoropentyl)-1H-indazol-3-yl (naphthalen-1-yl)methanone (THJ-2201);

the specific chemical designation (for purposes of this subsection only, the term "isomer" includes the optical, position and geometric isomers):

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

(2) 4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;

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

(4)(A) N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the acronym 25B-NBOMe.

(B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe).

(C) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)

(5) 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA;

(6) 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

(7) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

(8) 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-methylphenethylamine; paramethoxyamphetamine; PMA;

(9) 5-methoxy-3, 4-methylenedioxy-amphetamine;

(10) 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

(11) 3,4-methylenedioxy amphetamine;

(12) 3,4-methylenedioxymethamphetamine (MDMA);

(13) 3,4-methylenedioxy-N-ethylamphetamine (also known as – ethyl-alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

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

(15) 3,4,5-trimethoxy amphetamine;

(16) 5-methoxy-N, N-dimethyltryptamine (5-MeO-DMT);

(17) Alpha-methyltryptamine (other name: AMT);

(18) Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;

(19) Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;

(20) Dimethyltryptamine; some trade or other names: DMT;

(21) 5-Methoxy-N, N-diisopropyltryptamine (5-MeO-DIPT);

(22) 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;

(23) Lysergic acid diethylamide;

(24) Marijuana;

(25) Mescaline;

(26) 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; Synhexyl;

(27) 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 such plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of such plant, its seeds or extracts;

(28) N-ethyl-3-piperidyl benzilate;

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

(30) Psilocybin;

(31) Psilocyn;

(32) Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. and/or synthetic substances, immediate derivatives and their isomers with similar chemical structure and pharmacological activity such as the following:

delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;

delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;

delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;

(Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions covered).

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

(34) Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;

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

(36) 1[1-(2-thienyl)cyclohexyl]pyrroldine; some other names: TCPy.

(37) 4-methylmethcathinone (Mephedrone);

(38) 3,4-methylenedioxypyrovalerone (MDPV);

(39) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

(40) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

(41) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

(42) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

(43) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

(44) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

(45) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

(46) 2-(2,5-Dimethoxy-4-nitro-phenyl) ethanamine (2C-N);

(47) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);

(48) 3,4-Methylenedioxy-N-methylcathinone (Methylone);

(49) (2,5-dimethoxy-4-(n)-propyltghiophenethylamine (2C-T-7, itsoptical isomers, salts and salts of isomers

(50) 5-methoxy-N, N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT);

(51) Alpha-methyltryptamine (other name: AMT);

(52) 5-methoxy-N, N-diisopropyltryptamine (other name: 5-MeO-DIPT);

(53) Synthetic Cannabinoids as follows:

(A) 2-[(1R,3S)-3-hydroxycyclohexyl]-5- (2-methyloctan-2-yl) phenol) {also known as CP 47,497 and homologues};

(B) rel-2-[(1S,3R)-3-hydroxycyclohexyl] -5-(2-methylnonan-2-yl) phenol {also known as CP 47,497-C8 homolog};

(C) [(6a*R*)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7,10,10a-tetrahydrobenzo[c]chromen-1-ol)] {also known as HU-210};

(D) (dexanabinol);

(6a*S*,10a*S*)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzol[c]chromen-1-ol) {also known as HU-211};

(E) 1-Pentyl-3-(1-naphthoyl) indole {also known as JWH-018};

(F) 1-Butyl-3-(1-naphthoyl) indole {also known as JWH-073};

(G) (2-methyl-1-propyl-1H-indol-3-yl)-1-napthalenyl-methanone {also known as JWH-015};

(H) (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};

(I) [1-[2-(4-morpholinyl) ethyl] -1H-indol-3-yl]-1-naphthalenyl-methanone {also known as JWH-200};

(J) 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};

(K) 2-((1S,2S,5S)-5-hydroxy-2- (3-hydroxtpropyl)cyclohexyl) -5-(2-methyloctan-2-yl)phenol {also known as CP 55,940};

(L) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-122};

(M) (4-methyl-1-naphthalenyl) (1-pentyl-1H-indol-3-yl) -methanone {also known as JWH-398;

(N) (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};

(O) 1-(1-(2-cyclohexylethyl) -1H-indol-3-yl) -2-(2-methoxyphenyl) ethanone {also known as RCS-8};

(P) 1-pentyl-3-[1-(4-methoxynaphthoyl) indole (JWH-081);

(Q) 1-(5-fluoropentyl)-3-(1-naphthoyl) indole (AM2201); and

(R) 1-(5-fluoropentyl)-3-(2-iodobenzoyl) indole (AM694).

(54) Synthetic cannabinoids or any material, compound, mixture or preparation which contains any quantity of the following substances, including their analogues, congeners, homologues, isomers, salts and salts of analogues, congeners, homologues and isomers, as follows:

(A) CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-YL) phenol);

(B) HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10, 10A-tetrahydrobenzo[C] chromen-1-OL)];

(C) HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-YL)-6A,7,10,10 atetrahydrobenzo [C] chromen-1-OL);

(D) JWH-018, 1-pentyl-3-(1-naphthoyl) indole;

(E) JWH-019, 1-hexyl-3-(1-naphthoyl) indole;

(F) JWH-073, 1-butyl-3-(1-naphthoyl) indole;

(G) JWH-200, (1-(2-morpholin-4-ylethyl) indol-3-yl)- Naphthalen-1-ylmethanone;

(H) JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl) indole.

(55) Synthetic cannabinoids including any material, compound, mixture or preparation that is not listed as a controlled substance in Schedule I through V, is not a federal Food and Drug Administration approved drug or used within legitimate and approved medical research and which contains any quantity of the following substances, their salts, isomers, whether optical positional or geometric, analogues, homologues and salts of isomers, analogues and homologues, unless specifically exempted, whenever the existence of these salts, isomers, analogues, homologues and salts of isomers, analogues and homologues if possible within the specific chemical designation:

(A) Tetrahydrocannabinols meaning tetrahydrocannabinols which are naturally contained in a plant of the genus cannabis as well as synthetic equivalents of the substances contained in the plant or in the resinous extractives of cannabis or synthetic substances, derivatives and their isomers with analogous chemical structure and or pharmacological activity such as the following:

(i) DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.

(ii) DELTA-6 CIS OR trans tetrahydrocannabinol and their Optical isomers.

(iii) DELTA-3,4 CIS OR their trans tetrahydrocannabinol and their optical isomers.

(B) Naphthoyl indoles or any compound containing a 3-(-1- Napthoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include the following:

(i) JWH 015;

(ii) JWH 018;

(iii) JWH 019;

(iv) JWH 073;

(v) JWH 081;

(vi) JWH 122;

(vii) JWH 200;

(viii) JWH 210;

(ix) JWH 398;

(x) AM 2201;

(xi) WIN 55,212.

(56) Synthetic Phenethylamines (including their optical, positional, and geometric isomers, salts and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers):

(A) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-NBOMe);

(B) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-NBOMe);

(C) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-NBOMe);

(57) Synthetic Opioids (icluding their isomers, esters, ethers, salts and salts of isomers, esters and ethers):

(A) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);

(B) furanyl fentanyl;

(C) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-47700);

(D) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);

(E) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethylpiperidin-4-yl]-N-phenylpropionamide, also known as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide, (beta-hydroxythiofentanyl).

(58) Opioid Receptor Agonist (including its isomers, esters, ethers, salts, and salts of isomers, esters and ethers):

(A) AH-7921 (3,4-dichloro-N- (1dimethylamino)cyclohexylmethyl]benzamide).

(59) Naphylmethylindoles or any compound containing a 1hindol-3-yl-(1-naphthyl) methane structure with a substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 175 and JWH 184.

(60) Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure with substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the pyrrole ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 147 and JWH 307.

(61) Naphthylmethylindenes or any compound containing a Naphthylideneindene structure with substitution at the 3- Position of the indene ring whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall include, but not be limited to, JWH 176.

(62) Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include the following:

(A) RCS-8, SR-18 OR BTM-8;

(B) JWH 250;

(C) JWH 203;

(D) JWH 251;

(E) JWH 302.

(63) Cyclohexylphenols or any compound containing a 2-(3- hydroxycyclohexyl) phenol structure with a substitution at the 5-position of the phenolic ring whether or not substituted in the cyclohexyl ring to any extent. This shall include the following:

(A) CP 47,497 and its homologues and analogs;

(B) Cannabicyclohexanol;

(C) CP 55,940.

(64) Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with substitution at the nitrogren atom of the indole ring whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include the following:

(A) AM 694;

(B) Pravadoline WIN 48,098;

(C) RCS 4;

(D) AM 679.

(65) [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-napthalenymethanone. This shall include WIN 55,212-2.

(66) Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051 and JWH 133.

(67) Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with substitution at the nitrogen atom of the indole ring whether or not further substituted in the adamantoyl ring system to any extent. This shall include AM1248.

(68) Tetramethylcyclopropylindoles or any compound containing A 3-tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring

whether or not further substituted in the indole ring to any extent and whether or not substituted in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.

(69) N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.

(70) Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and V, not federal Food and Drug Administration approved drug or used within legitimate, approved medical research. Since nomenclature of these substances is not internationally standardized, any immediate precursor or immediate derivative of these substances shall be covered.

(71) Tryptamines:

(A) 5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MiPT)

(B) 4-hydroxy-N, N-diisopropyltryptamine (4-HO-DiPT)

(C) 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)

(D) 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)

(E) 4-acetoxy-N, N-diisopropyltryptamine (4-AcO-DiPT)

(F) 5-methoxy-α-methyltryptamine (5-MeO-AMT)

(G) 4-methoxy-N, N-Dimethyltryptamine (4-MeO-DMT)

(H) 4-hydroxy Diethyltryptamine (4-HO-DET)

(I) 5- methoxy- N, N- diallyltryptamine (5-MeO-DALT)

(J) 4-acetoxy-N, N-Dimethyltryptamine (4-AcO DMT)

(K) 4-hydroxy Diethyltryptamine (4-HO-DET)

(72) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (AB-CHMINACA);

(73) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (AB-PINACA);

(74) [1-(5-fluoropentyl)-1H-indazol-3-yl (naphthalen-1-yl)methanone (THJ-2201);

(75) quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB-22; QUPIC);

(76) quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22);

(77) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (AB-FUBINACA);

(78) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (ADB-PINACA); and

(79) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (common names, MAB-CHMINACA and ADB-CHMINACA);

(e) *Depressants*. — 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.

(f) *Stimulants.* — 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) Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine;

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

(3) Fenethylline;

(4) Methcathinone, its immediate precursors and immediate derivatives, its salts, optical isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-

(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or mephedrone;3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone; methylcathinone; AL-464; AL-422; AL- 463 and UR1432;

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

(6) N-ethylamphetamine;

(7) N,N-dimethylamphetemine; also known as N,N-alpha- trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine.

(8) Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and salts of isomers.

(9) Substituted amphetamines:

(A) 2-Fluoroamphetamine

(B) 3-Fluoroamphetamine

(C) 4-Fluoroamphetamine

(D) 2-chloroamphetamine

(E) 3-chloroamphetamine

(F) 4-chloroamphetamine

(G) 2-Fluoromethamphetamine

(H) 3-Fluoromethamphetamine

(I) 4-Fluoromethamphetamine

(J) 4-chloromethamphetamine

(10) 4-methyl-N-ethylcathinone (4-MEC);

(11) 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP);

(12) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone);

(13) 2-(methylamino)-1-phenylpentan-1-one (pentedrone);

(14) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone);

(15) 4-fluoro-N-methylcathinone (4-FMC);

(16) 3-fluoro-N-methylcathinone (3-FMC);

(17) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); and

(18) Alpha-pyrrolidinobutiophenone (α-PBP).

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

(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical 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.

(3) N-benzylpiperazine, also known as BZP.

(h) The following controlled substances are included in Schedule I:

(1) Synthetic Cathinones or any compound, except bupropion or compounds listed under a different schedule, or compounds used within legitimate and approved medical research, structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic or fused polycyclic ring systems, whether or not the compound is further modified in any of the following ways:

(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.

(D) By inclusion of the 2-amino nitrogen atom in a cyclic structure.

(2) Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV

and V, not federal Food and Drug Administration approved drug or used within legitimate, approved medical research.

§60A-2-206. Schedule II.

(a) Schedule II consists of the drugs and other substances, by whatever official name, common or usual name, chemical name or brand name designated, listed in this section.

(b) *Substances, vegetable origin or chemical synthesis.* — Unless specifically excepted or unless listed in another schedule, any of the following substances whether produced directly or indirectly by extraction from substances of vegetable origin, or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis:

(1) Opium and opiate, and any salt, compound, derivative or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, naloxone and naltrexone, and their respective salts, but including the following:

(A) Raw opium;

(B) Opium extracts;

(C) Opium fluid;

(D) Powdered opium;

(E) Granulated opium;

(F) Tincture of opium;

(G) Codeine;

(H) Dihydroetorphine;

(I) Ethylmorphine;

(J) Etorphine hydrochloride;

(K) Hydrocodone;

(L) Hydromorphone;

(M) Metopon;

(N) Morphine;

(O) Oripavine;

(P) Oxycodone;

(Q) Oxymorphone; and

(R) Thebaine;

(2) Any salt, compound, derivative or preparation thereof which is chemically equivalent or identical with any of the substances referred to in subdivision (1) of this subsection, except that these substances shall not include the isoquinoline alkaloids of opium;

(3) Opium poppy and poppy straw;

(4) Coca leaves and any salt, compound, derivative or preparation of coca leaves (including cocaine and ecgonine and their salts, isomers, derivatives and salts of isomers and derivatives), and any salt, compound, derivative or preparation thereof which is chemically equivalent or identical with any of these substances, except that the substances shall not include decocainized coca leaves or extractions of coca leaves, which extractions do not contain cocaine or ecgonine;

(5) Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid or powder form which contains the phenanthrene alkaloids of the opium poppy).

(c) *Opiates*. — Unless specifically excepted or unless in another schedule, any of the following opiates, including its 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, dextrorphan and levopropoxyphene excepted:

(1) Alfentanil;

(2) Alphaprodine;

(3) Anileridine;

(4) Bezitramide;

(5) Bulk dextropropoxyphene (nondosage forms);

(6) Carfentanil;

(7) Dihydrocodeine;

(8) Diphenoxylate;

(9) Fentanyl;

(10) Isomethadone;

(11) Levo-alphacetylmethadol; some other names: levo-alpha-acetylmethadol, levomethadyl acetate, LAAM;

(12) Levomethorphan;

(13) Levorphanol;

(14) Metazocine;

(15) Methadone;

(16) Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane;

(17) Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic acid;

(18) Pethidine; (meperidine);

(19) Pethidine-Intermediate-A, 4-cyano-1-methyl-4- phenylpiperidine;

(20) Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;

(21) Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;

(22) Phenazocine;

(23) Piminodine;

(24) Racemethorphan;

(25) Racemorphan;

(26) Remifentanil;

(27) Sufentanil;

(28) Tapentadol;

(29) Thiafentanil (4-(methoxycarbonyl)-4-(N-phenmethoxyacetamido)-1-2-(thienyl)ethylpiperidine), including its isomers, esters, ethers, salts and salts of isomers, esters and ethers.

(d) *Stimulants.* — 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:

(1) Amphetamine, its salts, optical isomers and salts of its optical isomers;

(2) Methamphetamine, its salts, isomers and salts of its isomers;

(3) Methylphenidate;

(4) Phenmetrazine and its salts; and

(5) Lisdexamfetamine.

(e) *Depressants.* — 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) Amobarbital;

(2) Glutethimide;

(3) Pentobarbital;

(4) Phencyclidine;

(5) Secobarbital.

(f) *Hallucinogenic substances*:

Nabilone: [Another name for nabilone: (+-)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8, 10, 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9H-dibenzo [b,d] pyran-9-one].

(g) *Immediate precursors.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

(1) Immediate precursor to amphetamine and methamphetamine:

(A) Phenylacetone;

(B) Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;

(2) Immediate precursors to phencyclidine (PCP):

(A) 1-phenylcyclohexylamine; and

(B) 1-piperidinocyclohexanecarbonitrile (PCC).

(3) Immediate precursor to fentanyl:

4-anilino-N-phenethyl-4-piperidine (ANPP).

§60A-2-210. Schedule IV.

(a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.

(b) *Narcotic drugs.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:

(1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit;

(2) Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane).

(c) *Depressants.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation which contains any quantity of the following substances, 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) Alprazolam;

(2) Barbital;

(3) Bromazepam;

(4) Camazepam;

(5) Carisoprodol;

(6) Chloral betaine;

(7) Chloral hydrate;

(8) Chlordiazepoxide;

(9) Clobazam;

(10) Clonazepam;

(11) Clorazepate;

(12) Clotiazepam;

(13) Cloxazolam;

(14) Delorazepam;

(15) Diazepam;

(16) Dichloralphenazone;

(17) Estazolam;

(18) Ethchlorvynol;

(19) Ethinamate;

(20) Ethyl loflazepate;

(21) Fludiazepam;

(22) Flunitrazepam;

(23) Flurazepam;

(24) Fospropofol;

(25) Halazepam;

(26) Haloxazolam;

(27) Ketazolam;

(28) Loprazolam;

(29) Lorazepam;

(30) Lormetazepam;

(31) Mebutamate;

(32) Medazepam;

(33) Meprobamate;

(34) Methohexital;

(35) Methylphenobarbital (mephobarbital);

(36) Midazolam;

(37) Nimetazepam;

(38) Nitrazepam;

(39) Nordiazepam;

(40) Oxazepam;

(41) Oxazolam;

(42) Paraldehyde;

(43) Petrichloral;

(44) Phenobarbital;

(45) Pinazepam;

(46) Prazepam;

(47) Quazepam;

(48) Temazepam;

(49) Tetrazepam;

(50) Triazolam;

(51) Zaleplon;

(52) Zolpidem;

(53) Zopiclone'

(75) quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB-22; QUPIC);

(76) quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22);

(77) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (AB-FUBINACA);

(78) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (ADB-PINACA); and

(79) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (common names, MAB-CHMINACA and ADB-CHMINACA);

(e) *Depressants.* — 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.

(f) *Stimulants.* — 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) Aminorex; some other names: aminoxaphen; 2-amino-5- phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine;

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

(3) Fenethylline;

(4) Methcathinone, its immediate precursors and immediate derivatives, its salts, optical isomers and salts of optical isomers; some other names: (2-(methylamino)-propiophenone; alpha-

(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1- one; alpha—methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypyrovalerone and/or mephedrone;3,4-methylenedioxypyrovalerone (MPVD); ephedrone; N-methylcathinone; methylcathinone; AL-464; AL-422; AL- 463 and UR1432;

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

(6) N-ethylamphetamine;

(7) N,N-dimethylamphetemine; also known as N,N-alpha- trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine.

(8) Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and salts of isomers.

(9) Substituted amphetamines:

(A) 2-Fluoroamphetamine

(B) 3-Fluoroamphetamine

(C) 4-Fluoroamphetamine

(D) 2-chloroamphetamine

(E) 3-chloroamphetamine

(F) 4-chloroamphetamine

(G) 2-Fluoromethamphetamine

(H) 3-Fluoromethamphetamine

(I) 4-Fluoromethamphetamine

(J) 4-chloromethamphetamine

(10) 4-methyl-N-ethylcathinone (4-MEC);

(11) 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP);

(12) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone);

(13) 2-(methylamino)-1-phenylpentan-1-one (pentedrone);

(14) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone);

(15) 4-fluoro-N-methylcathinone (4-FMC);

(16) 3-fluoro-N-methylcathinone (3-FMC);

(17) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); and

(18) Alpha-pyrrolidinobutiophenone (α-PBP).

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

(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical 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.

(3) N-benzylpiperazine, also known as BZP.

(h) The following controlled substances are included in Schedule I:

(1) Synthetic Cathinones or any compound, except bupropion or compounds listed under a different schedule, or compounds used within legitimate and approved medical research, structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic or fused polycyclic ring systems, whether or not the compound is further modified in any of the following ways:

(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.

(D) By inclusion of the 2-amino nitrogen atom in a cyclic structure.

(2) Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV

and V, not federal Food and Drug Administration approved drug or used within legitimate, approved medical research.

§60A-2-206. Schedule II.

(a) Schedule II consists of the drugs and other substances, by whatever official name, common or usual name, chemical name or brand name designated, listed in this section.

(b) *Substances, vegetable origin or chemical synthesis.* — Unless specifically excepted or unless listed in another schedule, any of the following substances whether produced directly or indirectly by extraction from substances of vegetable origin, or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis:

(1) Opium and opiate, and any salt, compound, derivative or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, naloxone and naltrexone, and their respective salts, but including the following:

(A) Raw opium;

(B) Opium extracts;

(C) Opium fluid;

(D) Powdered opium;

(E) Granulated opium;

(F) Tincture of opium;

(G) Codeine;

(H) Dihydroetorphine;

(I) Ethylmorphine;

(J) Etorphine hydrochloride;

(K) Hydrocodone;

(L) Hydromorphone;

(M) Metopon;

(N) Morphine;

(O) Oripavine;

(P) Oxycodone;

(Q) Oxymorphone; and

(R) Thebaine;

(2) Any salt, compound, derivative or preparation thereof which is chemically equivalent or identical with any of the substances referred to in subdivision (1) of this subsection, except that these substances shall not include the isoquinoline alkaloids of opium;

(3) Opium poppy and poppy straw;

(4) Coca leaves and any salt, compound, derivative or preparation of coca leaves (including cocaine and ecgonine and their salts, isomers, derivatives and salts of isomers and derivatives), and any salt, compound, derivative or preparation thereof which is chemically equivalent or identical with any of these substances, except that the substances shall not include decocainized coca leaves or extractions of coca leaves, which extractions do not contain cocaine or ecgonine;

(5) Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid or powder form which contains the phenanthrene alkaloids of the opium poppy).

(c) *Opiates.* — Unless specifically excepted or unless in another schedule, any of the following opiates, including its 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, dextrorphan and levopropoxyphene excepted:

(1) Alfentanil;

(2) Alphaprodine;

(3) Anileridine;

(4) Bezitramide;

(5) Bulk dextropropoxyphene (nondosage forms);

(6) Carfentanil;

(7) Dihydrocodeine;

(8) Diphenoxylate;

(9) Fentanyl;

(10) Isomethadone;

(11) Levo-alphacetylmethadol; some other names: levo-alpha-acetylmethadol, levomethadyl acetate, LAAM;

(12) Levomethorphan;

(13) Levorphanol;

(14) Metazocine;

(15) Methadone;

(16) Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane;

(17) Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic acid;

(18) Pethidine; (meperidine);

(19) Pethidine-Intermediate-A, 4-cyano-1-methyl-4- phenylpiperidine;

(20) Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;

(21) Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;

(22) Phenazocine;

(23) Piminodine;

(24) Racemethorphan;

(25) Racemorphan;

(26) Remifentanil;

(27) Sufentanil;

(28) Tapentadol;

(29) Thiafentanil (4-(methoxycarbonyl)-4-(N-phenmethoxyacetamido)-1-2-(thienyl)ethylpiperidine), including its isomers, esters, ethers, salts and salts of isomers, esters and ethers.

(d) *Stimulants.* — 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:

(1) Amphetamine, its salts, optical isomers and salts of its optical isomers;

(2) Methamphetamine, its salts, isomers and salts of its isomers;

(3) Methylphenidate;

(4) Phenmetrazine and its salts; and

(5) Lisdexamfetamine.

(e) *Depressants.* — 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) Amobarbital;

(2) Glutethimide;

(3) Pentobarbital;

(4) Phencyclidine;

(5) Secobarbital.

(f) *Hallucinogenic substances*:

Nabilone: [Another name for nabilone: (+-)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8, 10, 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9H-dibenzo [b,d] pyran-9-one].

(g) *Immediate precursors.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

(1) Immediate precursor to amphetamine and methamphetamine:

(A) Phenylacetone;

(B) Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;

(2) Immediate precursors to phencyclidine (PCP):

(A) 1-phenylcyclohexylamine; and

(B) 1-piperidinocyclohexanecarbonitrile (PCC).

(3) Immediate precursor to fentanyl:

4-anilino-N-phenethyl-4-piperidine (ANPP).

§60A-2-210. Schedule IV.

(a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.

(b) *Narcotic drugs.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:

(1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit;

(2) Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane).

(c) *Depressants.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation which contains any quantity of the following substances, 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) Alprazolam;

(2) Barbital;

(3) Bromazepam;

(4) Camazepam;

(5) Carisoprodol;

(6) Chloral betaine;

(7) Chloral hydrate;

(8) Chlordiazepoxide;

(9) Clobazam;

(10) Clonazepam;

(11) Clorazepate;

(12) Clotiazepam;

(13) Cloxazolam;

(14) Delorazepam;

(15) Diazepam;

(16) Dichloralphenazone;

(17) Estazolam;

(18) Ethchlorvynol;

(19) Ethinamate;

(20) Ethyl loflazepate;

(21) Fludiazepam;

(22) Flunitrazepam;

(23) Flurazepam;

(24) Fospropofol;

(25) Halazepam;

(26) Haloxazolam;

(27) Ketazolam;

(28) Loprazolam;

(29) Lorazepam;

(30) Lormetazepam;

(31) Mebutamate;

(32) Medazepam;

(33) Meprobamate;

(34) Methohexital;

(35) Methylphenobarbital (mephobarbital);

(36) Midazolam;

(37) Nimetazepam;

(38) Nitrazepam;

(39) Nordiazepam;

(40) Oxazepam;

(41) Oxazolam;

(42) Paraldehyde;

(43) Petrichloral;

(44) Phenobarbital;

(45) Pinazepam;

(46) Prazepam;

(47) Quazepam;

(48) Temazepam;

(49) Tetrazepam;

(50) Triazolam;

(51) Zaleplon;

(52) Zolpidem;

(53) Zopiclone'

(54) Suvorexant ([(7R)-4-(5-chloro-1,3-benzoxazol-2-yl)-7-methyl-1,4-diazepan-1-yl] [5-methyl-2-(2H-1,2,3-triazol-2-yl)phenyl]methanone).

(d) Any material, compound, mixture or preparation which contains any quantity of the following substance, including its salts, isomers (whether optical, position or geometric) and salts of such isomers whenever the existence of such salts, isomers and salts of isomers is possible: Fenfluramine and Dexfenfluramine.

(e) *Stimulants.* — 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) Cathine ((+)-norpseudoephedrine);

(2) Diethylpropion;

(3) Fencamfamin;

(4) Fenproporex;

(5) Mazindol;

(6) Mefenorex;

(7) Modafinil;

(8) Pemoline (including organometallic complexes and chelates thereof);

(9) Phentermine;

(10) Pipradrol;

(11) Sibutramine;

(12) SPA ((-)-1-dimethylamino-1,2-diphenylethane);

(13) Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropyl [(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid);

(f) *Other substances.* — Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation which contains any quantity of the following substances, including its salts:

(1) Pentazocine;

(2) Butorphanol;

(3) Tramadol (2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol).

Amyl nitrite, butyl nitrite, isobutyl nitrite and the other organic nitrites are controlled substances and no product containing these compounds as a significant component shall be possessed, bought or sold other than pursuant to a bona fide prescription or for industrial or manufacturing purposes.

§60A-2-212. Schedule V.

(a) Schedule V shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.

(b) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture or preparation containing any of the following narcotic drugs or their salts calculated as the free anhydrous base or alkaloid in limited quantities as set forth below, which shall include one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:

(1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;

(2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams;

(3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams;

(4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit;

(5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams;

(6) Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.

(c) *Stimulants.* — Unless specifically exempted or excluded 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) Pyrovalerone.

(d) Any compound, mixture or preparation containing as its single active ingredient ephedrine, pseudoephedrine or phenylpropanolamine, their salts or optical isomers, or salts of optical isomers except products which are for pediatric use primarily intended for administration to children under the age of twelve: *Provided,* That neither the offenses set forth in section four hundred one, article four of this chapter, nor the penalties therein, shall be applicable to ephedrine, pseudoephedrine or phenylpropanolamine which shall be subject to the provisions of article ten of this chapter.

(e) *Depressants.* — Unless specifically exempted or excluded 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:

(1) Ezogabine [N-[2-amino-4-94-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester];

(2) Lacosamide [(R)-2-acetoamido- *N* -benzyl-3-methoxy-propionamide];

(3) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid]; and

(4) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl] butanamide) (also referred to as BRV; UCB-34714; Briviact), including its salts.

The Joint Committee on Enrolled Bills hereby certifies that the foregoing bill is correctly enrolled.

Vice- Chairman, House Committee

Member ~~Chairman,~~ Senate Committee

Originating in the House.

In effect ninety days from passage.

Clerk of the House of Delegates

Clerk of the Senate

Speaker of the House of Delegates

President of the Senate

FILED
2017 APR 26 P 8:19
OFFICE WEST VIRGINIA
SECRETARY OF STATE

The within is approved this the 26th

day of April, 2017.

Governor

PRESENTED TO THE GOVERNOR

APR 21 2017

Time 3:57 pm