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STATE OF HAWAII | KA MOKUʻĀINA O HAWAIʻI
DEPARTMENT OF PUBLIC SAFETY
KA ʻOIHANA HOʻOPALEKANA LEHULEHU
1177 Alakea Street
Honolulu, Hawaii 96813

No. _____

TESTIMONY ON HOUSE BILL 1097, HOUSE DRAFT (HD) 1
RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT
By
Tommy Johnson, Director

House Committee on Judiciary and Hawaiian Affairs
Representative David A.Tarnas, Chair
Representative Gregg Takayama, Vice Chair

Thursday, February 23, 2023; 2:00 p.m.
Via Videoconference

Chair Tarnas, Vice Chair Takayama, and Members of the Committee:

The Department of Public Safety (PSD) **supports** House Bill (HB) 1097, HD 1, which updates chapter 329 of the Hawaii Revised Statutes (HRS), to incorporate amendments made to the federal Controlled Substances Act that were temporarily permitted in Hawaii by the temporary designation of several new controlled substances by PSD in 2022. Under section 329-11(d) and (e), HRS, PSD's temporary designation of new controlled substances shall be nullified if the next regular session of the State Legislature has not made the corresponding changes to state law.

Sections 329-11(d) and (e), HRS, provide that if a substance is added, deleted, or rescheduled under federal law, or by an emergency scheduling action taken by PSD, then PSD shall recommend to the Legislature to make the corresponding changes in Hawaii law. In 2022, PSD temporarily designated many new controlled substances, and those substances are listed in HB 1097, HD 1.

HB 1097, also amends chapter 329, HRS, to mirror recent changes to the federal Controlled Substances Act, thereby eliminating differences between federal and state law and avoiding nullification of the controlled substances that were temporarily designated in 2022.

PSD respectfully requests that the committee amend HB 1097, HD 1, because drafting errors in the original proposal were discovered requiring changes to various parts of HB 1097, HD 1.

Requested Amendments to HD1:

SECTION 1. Section 329-14, Hawaii Revised Statutes, is amended as follows:

1. By amending subsection (b) to read as follows:

"(b) Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

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

(2) Acetylmethadol;

(3) Acetyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide), its optical, positional, and geometric isomers, salts, and salts of isomers;

(4) Acryl fentanyl [N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide] (Other name: acryloylfentanyl);

(5) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)cyclohexylmethyl]benzamide);

(6) Allylprodine;

(7) Alphacetylmethadol (except levo-alphacetylmethadol, levomethadyl acetate, or LAAM);

(8) Alphameprodine;

(9) Alphamethadol;

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

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

(12) Benzethidine;

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

(14) Betacetylmethadol;

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

(16) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide);

(17) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide);

(18) Betameprodine;

(19) Betamethadol;

(20) Beta-methyl fentanyl (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide) (Other name: [beta]-methyl fentanyl);

(21) Beta'-phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide) (Other names: [beta]'-phenyl fentanyl; 3-phenylpropanoyl fentanyl);

(22) Betaprodine;

(23) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide);

(24) Clonitazene;

(25) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);

(26) Dextromoramide;

(27) Diampromide;

(28) Diethylthiambutene;

(29) Difenoxin;

(30) Dimenoxadol;

(31) Dimepheptanol;

(32) Dimethylthiambutene;

(33) Dioxaphetyl butyrate;

(34) Dipipanone;

(35) Ethylmethylthiambutene;

(36) Etonitazene;

(37) Etoxeridine;

(38) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-yl)(phenyl)carbamate);

(39) 4-fluoroisobutyryl fentanyl [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide] (Other name: para-fluoroisobutyryl fentanyl);

(40) 2'-fluoro ortho-fluorofentanyl (N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide) (Other name: 2'-fluoro 2-fluorofentanyl);

(41) Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide);

(42) Furethidine;

(43) Hydroxypethidine;

(44) Ketobemidone;

(45) Levomoramide;

(46) Levophenacylmorphan;

(47) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

(48) 4'-methyl acetyl fentanyl (N-(1-(4-methylphenethyl)piperidin-4-yl)-N-phenylacetamide);

(49) 3-methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

(50) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

(51) Morpheridine;

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

(53) Noracymethadol;

(54) Norlevorphanol;

(55) Normethadone;

(56) Norpipanone;

(57) Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide];

(58) Ortho-fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide);

(59) Ortho-fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) (Other name: 2-fluorobutyryl fentanyl);

(60) Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide) (Other name: 2-fluorofentanyl);

(61) Ortho-fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);

(62) Ortho-methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide) (Other name: 2-methyl acetylfentanyl);

(63) Ortho-methyl methoxyacetyl fentanyl (2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide) (Other name: 2-methyl methoxyacetyl fentanyl);

(64) Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);

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

(66) Para-fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide);

(67) Para-methylfentanyl (N-(4-methylphenyl)-N(1-phenethylpiperidin-4-yl)propionamide) (Other name: 4-methylfentanyl);

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

(69) Phenadoxone;

(70) Phenampromide;

(71) Phenomorphan;

(72) Phenoperidine;

(73) Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide) (Other name: benzoyl fentanyl);

(74) Piritramide;

(75) Proheptazine;

(76) Properidine;

(77) Propiram;

(78) Racemoramide;

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

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

(81) Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide) (Other names: 2-thiofuranyl fentanyl; thiophene fentanyl);

(82) Tilidine;

(83) Trimeperidine; [and]

(84) U-47700 (3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide)[.].”

2. By amending subsection (d) to read as follows:

“(d) Any material, compound, mixture, or preparation that contains any quantity of the following hallucinogenic substances, their salts, isomers, and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Alpha-ethyltryptamine (AET);

(2) 2,5-dimethoxy-4-ethylamphetamine (DOET);

(3) 2,5-dimethoxyamphetamine (2,5-DMA);

(4) 3,4-methylenedioxy amphetamine;

(5) 3,4-methylenedioxymethamphetamine (MDMA);

(6) N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-MDA);

(7) 3,4-methylenedioxy-N-ethylamphetamine (MDE);

(8) 5-methoxy-3,4-methylenedioxy-amphetamine;

(9) 4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);

(10) 4-Bromo-2,5-dimethoxyphenethylamine (Nexus);

(11) 3,4,5-trimethoxy amphetamine;

(12) Bufotenine;

(13) 4-methoxyamphetamine (PMA);

(14) Diethyltryptamine;

(15) Dimethyltryptamine;

(16) 4-methyl-2,5-dimethoxy-amphetamine;

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

(18) Ibogaine;

(19) Lysergic acid diethylamide;

(20) Marijuana;

(21) Parahexyl;

(22) Mescaline;

(23) Peyote;

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

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

(26) Psilocybin;

(27) Psilocyn;

(28) 1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);

(29) Ethylamine analog of phencyclidine (PCE);

(30) Pyrrolidine analog of phencyclidine (PCPy, PHP);

(31) Thiophene analog of phencyclidine (TPCP; TCP);

(32) Gamma-butyrolactone, including butyrolactone; butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone dihydro; dihydro-2(3H)furanone; tetrahydro-2-furanone; 1,2-butanolide; 1,4-butanolide; 4-butanolide; gamma-hydroxybutyric acid lactone; 3-hydroxybutyric acid lactone and 4-hydroxybutanoic acid lactone with Chemical Abstract Service number 96-48-0 when any such substance is intended for human ingestion;

(33) 1,4 butanediol, including butanediol; butane-1,4-diol; 1,4- butylenes glycol; butylene glycol; 1,4-dihydroxybutane; 1,4- tetramethylene glycol; tetramethylene glycol; tetramethylene 1,4- diol with Chemical Abstract Service number 110-63-4 when any such substance is intended for human ingestion;

(34) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7), its optical isomers, salts, and salts of isomers;

(35) N-benzylpiperazine (BZP; 1-benzylpiperazine) its optical isomers, salts, and salts of isomers;

(36) 1-(3-trifluoromethylphenyl)piperazine (TFMPP), its optical isomers, salts, and salts of isomers;

(37) Alpha-methyltryptamine (AMT), its isomers, salts, and salts of isomers;

(38) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT), its isomers, salts, and salts of isomers;

(39) Salvia divinorum;

(40) Salvinorin A;

(41) Divinorin A;

(42) 5-Methoxy-N,N-Dimethyltryptamine (5-MeO-DIPT) (some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT);

(43) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

(44) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

(45) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

(46) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

(47) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

(48) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

(49) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

(50) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);

(51) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);

(52) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5);

(53) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82); ~~and~~

(54) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);

(55) 2-(ethylamino)-2-(3-methoxylphenyl)cyclohexan-1-one(methoxetamine, MXE), 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;

(56) N-ethylhexedrone (Other names: aethylaminohexanophenone, 2-(ethylamino)-1-phenylhexan-s1-one);

(57) alpha-pyrrolidinohexanophenone (Other names: a-PHP, apyrrolidinohexanophenone, 1-phenyl-2-(pyrrolidine-1-yl)hexan-1-one);

(58) 4-methyl-alphaethylaminopentiophenone (Other names: 4-MEAP, 2-(ethylamino)-1-(4-methylphyl)pentan-1-one);

(59) 4'-methyl-alphapyrrolidinohexiophenone (Other names: MPHP, 4'-methyl-alphapyrrolidinohexanophenone, 1-(4-methylphenyl)-2-(pyrrolidine-1-yl)hexan1-one);

(60) Alpha-pyrrolidinoheptaphenone (Other names: PV8, 1-phenyl-2-(pyrrolidine-1-yl)heptan-1-one); and;

(61) 4'chloro-alphapyrrolidinovalerophenone (Other names: 4-chloro-a-PVP, 4'-chloro-apyrrolidinopentiophenone, 1-(4-chlorophenyl)-2-(pyrrolidine-1-yl)pentan-1-one) including their optical, positional, and geometric isomers, salts, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."

3. By amending subsection (e) to read as follows:

"(e) Depressants. Unless specifically excepted, the schedule shall include any material, compound, mixture, or preparation which contains any quantity of the substance:

(1) Mecloqualone;

(2) Methaqualone;

(3) Etizolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible); [or]

(4) Flualprazolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible)[.];

~~(5) N-ethylhexedrone (Other names: aethylaminohexanophenone, 2-(ethylamino)-1-phenylhexan-s1-one);~~

~~(6) alpha-pyrrolidinohexanophenone (Other names: a-PHP, apyrrolidinohexanophenone, 1-phenyl-2-(pyrrolidine-1-yl)hexan-1-one);~~

~~(7) 4-methyl-alphaethylaminopentiophenone (Other names: 4-MEAP, 2-(ethylamino)-1-(4-methylphyl)pentan-1-one);~~

~~(8) 4'-methyl-alphapyrrolidinohexiophenone (Other names: MPHP, 4'-methyl-alphapyrrolidinohexanophenone, 1-(4-methylphenyl)-2-(pyrrolidine-1-yl)hexan1-one);~~

~~(9) Alpha-pyrrolidinoheptaphenone (Other names: PV8, 1-phenyl-2-(pyrrolidine-1-yl)heptan-1-one); and~~

~~(10) 4'chloro-alphapyrrolidinovalerophenone (Other names: 4-chloro-a-PVP, 4'-chloro-apyrrolidinopentiophenone, 1-(4-chlorophenyl)-2-(pyrrolidine-1-yl)pentan-1-one) including their optical, positional, and geometric isomers, salts, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."~~

4. By amending subsection (f) to read as follows:

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

(1) Aminorex;

(2) Cathinone;

(3) 4,4'-dimethylaminorex (common name: 4,4'-DMAR);

(4) Fenethylline;

(5) Methcathinone;

(6) 4-methylaminorex;

(7) N-ethylamphetamine;

(8) N,N-dimethylamphetamine;

(9) Substituted cathinones, any compound, except bupropion or compounds listed under a different schedule, structurally derived from 2-aminopropan-1-one by substitution at the 1-position with either phenyl, naphthyl, or thiophene ring systems, whether or not the compound is further modified 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; or

(C) 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.

Some other trade names: Mephedrone (2-methylamino-1-p-tolylpropan-1-one), also known as 4-methylmethcathinone (4-MMC), methylephedrone or MMCAT; Methylenedioxypyrovalerone (MDPV, MDPK); methylone or 3,4-methylenedioxymethcathinone; and 1-(benzo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one, monohydrochloride, also known as Ethylone, bk-MDEA hydrochloride, MDEC; 3,4-Methylenedioxy-N-ethylcathinone; bk-Methylenedioxyethylamphetamine, 4-methyl-N-ethylcathinone (4-MEC); 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP); alpha-pyrrolidinopentiophenone ([alpha]-PVP); 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-1-one (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone, bk-MBDP); 4-fluoro-N-methylcathinone (4-FMC, flephedrone); 3-fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-pyrrolidinobutiophenone ([alpha]-PBP) and their optical, positional, and geometric isomers, salts and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers is possible; and

(10) 1-(4-methoxyphenyl)-N-methylpropan-2-amine (Other names: para-methoxymethamphetamine; PMMA)[.];

~~(11) Ethyl 2-(1-(5-fluoropentyl)-1Hindazole-3-carboxamido)-3,3-demethylbutanoate (Other name: 5FEDMB-PINACA);~~

~~(12) Methyl 2-(1-(5-fluropentyl)-1Hindole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5FMDMB-PICA; 5F-MDMB-2201);~~

~~(13) N-(Adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-fluorobenzyl));~~

~~(14) 1-(5-Fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (Other names: 5F-CUMYLPINACA; SGT-25); and~~

~~(15) (1-(4-Fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144) including their salts, isomers, and salts of isomers whenever the~~

existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."

5. By amending subsection (g) to read as follows:

(g) Any of the following cannabinoids, their salts, isomers, and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Tetrahydrocannabinols; meaning tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. or synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following: Delta 1 cis or trans tetrahydrocannabinol, and their optical isomers; Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers; and 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, are covered);

(2) Naphthoylindoles; meaning any compound containing a 3-(1-naphthoyl)indole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 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 naphthyl ring to any extent;

(3) Naphthylmethylindoles; meaning any compound containing a 1H-indol-3-yl-(1-naphthyl) methane structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 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 naphthyl ring to any extent;

(4) Naphthoylpyrroles; meaning any compound containing a 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom of the pyrrole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the pyrrole ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(5) Naphthylmethylindenes; meaning any compound containing a naphthylideneindene structure with substitution at the 3-position of the indene ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the indene ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(6) Phenylacetylindoles; meaning any compound containing a 3-phenylacetylindole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the indole ring to any extent, whether or not substituted in the phenyl ring to any extent;

(7) Cyclohexylphenols; meaning any compound containing a 2-(3-hydroxycyclohexyl) phenol structure with substitution at the 5-position of the phenolic ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not substituted in the cyclohexyl ring to any extent;

(8) Benzoylindoles; meaning any compound containing a 3-(benzoyl) indole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 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;

(9) [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl) pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-naphthalenylmethanone (another trade name is WIN 55,212-2);

(10) (6a,10a)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol (Other trade names are: HU-210/HU-211);

(11) Tetramethylcyclopropanoylindoles; meaning any compound containing a 3-tetramethylcyclopropanoylindole structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the tetramethylcyclopropyl ring to any extent;

(12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: APINACA, AKB48);

(13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: PB-22; QUPIC);

(14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-carboxylate, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 5-fluoro-PB-22; 5F-PB-22);

(15) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-FUBINACA);

(16) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: ADB-PINACA);

(17) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-CHMINACA);

(18) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: AB-PINACA);

(19) [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl)methanone, and geometric isomers, salts, and salts of isomers (Other names: THJ-2201);

(20) Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate, and geometric isomers, salts, and salts of isomers (Other names: FUB-AMB, Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);

(21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, and geometric isomers, salts, and salts of isomers (Other names: 5-fluoro-AMB, 5-fluoro-AMP);

(22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: AKB48 N-(5-fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl analog, 5F-APINACA);

(23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and geometric isomers, salts, and salts of isomers (Other names: STS-135, 5F-APICA; 5-fluoro-APICA);

(24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate, and geometric isomers, salts, and salts of isomers (Other names: NM2201; CBL2201);

(25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: MAB-CHMINACA and ADB-CHMINACA);

(26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (Other names: 5F-ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical, positional, and geometric isomers, salts, and salts of isomers;

(27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA; CUMYL-4CN-BINACA);

(28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (Other name: 5F-AB-PINACA);

(29) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoate (Other names: MMB-CHMICA; AMB-CHMICA);

(30) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-CUMYL-P7AICA); ~~and~~

(31) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate (MDMB-4en-PINACA~~.~~;

(32) Ethyl 2-(1-(5-fluoropentyl)-1Hindazole-3-carboxamido)-3,3-demethylbutanoate (Other name: 5FEDMB-PINACA);

(33) Methyl 2-(1-(5-fluropentyl)-1Hindole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5FMDMB-PICA; 5F-MDMB-2201);

(34) N-(Adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-fluorobenzyl));

(35) 1-(5-Fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (Other names: 5F-CUMYLPINACA; SGT-25); and

(36) (1-(4-Fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144) including their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."

SECTION 2. Section 329-20, Hawaii Revised Statutes, is amended by amending subsection (d) to read as follows:

1. By amending subsection (b) to read as follows:

"(b) Depressants. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, esters, ethers, and salts of isomers, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation, that has a degree of danger or probable danger associated with a depressant effect on the central nervous system:

(1) Alprazolam;

(2) Barbital;

(3) Brexanolone;

(4) Bromazepam;

(5) Butorphanol;

(6) Camazepam;

(7) Carisoprodol;

(8) Chloral betaine;

(9) Chloral hydrate;

(10) Chlordiazepoxide;

(11) Clobazam;

(12) Clonazepam;

(13) Clorazepate;

(14) Clotiazepam;

(15) Cloxazolam;

(16) Delorazepam;

(17) Diazepam;

(18) Dichloralphenazone (Midrin);

(19) Estazolam;

(20) Ethchlorvynol;

(21) Ethinamate;

(22) Ethyl loflazepate;

(23) Fludiazepam;

(24) Flunitrazepam;

(25) Flurazepam;

(26) Fospropofol (Lusedra);

(27) Halazepam;

(28) Haloxazolam;

(29) Ketazolam;

(30) Lemborexant ((1R,2S)-2-[(2,4-dimethylpyrimidin-5-yl)oxymethyl]-2-(3-fluorophenyl)-N-(5-fluoropyridin-2-yl)cyclopropane-1-carboxamide);

(31) Loprazolam;

(32) Lorazepam;

(33) Lormetazepam;

(34) Mebutamate;

(35) Medazepam;

(36) Meprobamate;

(37) Methohexital;

(38) Methylphenobarbital (mephorbarbital);

(39) Midazolam;

(40) Nimetazepam;

(41) Nitrazepam;

(42) Nordiazepam;

(43) Oxazepam;

(44) Oxazolam;

(45) Paraldehyde;

(46) Petrichloral;

(47) Phenobarbital;

(48) Pinazepam;

(49) Prazepam;

(50) Quazepam;

(51) Remimazolam;

(52) Suvorexant;

(53) Temazepam;

(54) Tetrazepam;

(55) Triazolam;

(56) Zaleplon;

(57) Zolpidem~~.~~; and

(58) Zopiclone (Lunesta)~~.~~; and

(59) Daridorexant and its salts."

2. By amending subsection (d) to read as follows:

"(d) Stimulants. Unless listed in another schedule, any material, compound, mixture, or preparation [which] that contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Cathine ((+)-norpseudoephedrine);

(2) Diethylpropion;

(3) Fencamfamin;

(4) Fenproporex;

(5) Lorcaserin;

(6) Mazindol;

(7) Mefenorex;

(8) Modafinil;

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

(10) Phentermine;

(11) Pipradrol;

(12) Serdexmethylphenidate;

(13) Sibutramine;

(14) Solriamfetol; [and]

(15) SPA (1-dimethylamino-1,2-dephenylethane, lefetamine)[.];

~~(16)~~ ~~Serdexmethylphenidate, including its salts, isomers, and salts of isomers; and~~

~~(17)~~ ~~Daridorexant and its salts.~~"

Thank you for the opportunity to testify in this measure.

JOSH GREEN
GOVERNOR

JORDAN LOWE
DIRECTOR

MICHAEL S. VINCENT
Deputy Director
Administration

Vacant
Deputy Director
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No.

TESTIMONY ON HOUSE BILL 1097, HOUSE DRAFT (HD) 1
RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT
Before the Senate Committee on Judiciary and Hawaiian Affairs
Thursday, February 23, 2023; 2:00 p.m.
State Capitol Conference Room 325, Via Videoconference
WRITTEN TESTIMONY ONLY

Chair Tarnas, Vice Chair Takayama, and Members of the Committee:

The Department of Law Enforcement (DLE) **supports** House Bill (HB) 1097, HD 1, which updates chapter 329 of the Hawaii Revised Statutes (HRS), to incorporate amendments made to the federal Controlled Substances Act that were temporarily permitted in Hawaii by the temporary designation of several new controlled substances by PSD in 2022. Under section 329-11(d) and (e), HRS, DLE's temporary designation of new controlled substances shall be nullified if the next regular session of the State Legislature has not made the corresponding changes to state law.

Sections 329-11(d) and (e), HRS, provide that if a substance is added, deleted, or rescheduled under federal law, or by an emergency scheduling action taken by DLE, then DLE shall recommend to the Legislature to make the corresponding changes in Hawaii law. In 2022, PSD temporarily designated many new controlled substances, and those substances are listed in HB 1097, HD 1.

HB 1097, also amends chapter 329, HRS, to mirror recent changes to the federal Controlled Substances Act, thereby eliminating differences between federal and state law and avoiding nullification of the controlled substances that were temporarily designated in 2022.

DLE respectfully requests that the committee amend HB 1097, HD 1, because drafting errors in the original proposal were discovered requiring changes to various parts of HB 1097, HD 1.

Requested Amendments to HD1:

SECTION 1. Section 329-14, Hawaii Revised Statutes, is amended as follows:

1. By amending subsection (b) to read as follows:

"(b) Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

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

(2) Acetylmethadol;

(3) Acetyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide), its optical, positional, and geometric isomers, salts, and salts of isomers;

(4) Acryl fentanyl [N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide] (Other name: acryloylfentanyl);

(5) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)cyclohexylmethyl]benzamide);

(6) Allylprodine;

(7) Alphacetylmethadol (except levo-alphacetylmethadol, levomethadyl acetate, or LAAM);

(8) Alphameprodine;

(9) Alphamethadol;

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

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

(12) Benzethidine;

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

(14) Betacetylmethadol;

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

(16) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide);

(17) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide);

(18) Betameprodine;

(19) Betamethadol;

(20) Beta-methyl fentanyl (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide) (Other name: [beta]-methyl fentanyl);

(21) Beta'-phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide) (Other names: [beta]'-phenyl fentanyl; 3-phenylpropanoyl fentanyl);

(22) Betaprodine;

(23) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide);

(24) Clonitazene;

(25) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);

(26) Dextromoramide;

(27) Diampromide;

(28) Diethylthiambutene;

(29) Difenoxin;

(30) Dimenoxadol;

(31) Dimepheptanol;

(32) Dimethylthiambutene;

(33) Dioxaphetyl butyrate;

(34) Dipipanone;

(35) Ethylmethylthiambutene;

(36) Etonitazene;

(37) Etoxeridine;

(38) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-yl)(phenyl)carbamate);

(39) 4-fluoroisobutyryl fentanyl [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide] (Other name: para-fluoroisobutyryl fentanyl);

(40) 2'-fluoro ortho-fluorofentanyl (N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide) (Other name: 2'-fluoro 2-fluorofentanyl);

(41) Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide);

(42) Furethidine;

(43) Hydroxypethidine;

(44) Ketobemidone;

(45) Levomoramide;

(46) Levophenacylmorphan;

(47) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

(48) 4'-methyl acetyl fentanyl (N-(1-(4-methylphenethyl)piperidin-4-yl)-N-phenylacetamide);

(49) 3-methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

(50) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

(51) Morpheridine;

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

(53) Noracymethadol;

(54) Norlevorphanol;

(55) Normethadone;

(56) Norpipanone;

(57) Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide];

(58) Ortho-fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide);

(59) Ortho-fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) (Other name: 2-fluorobutyryl fentanyl);

(60) Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide) (Other name: 2-fluorofentanyl);

(61) Ortho-fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);

(62) Ortho-methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide) (Other name: 2-methyl acetylfentanyl);

(63) Ortho-methyl methoxyacetyl fentanyl (2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide) (Other name: 2-methyl methoxyacetyl fentanyl);

(64) Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);

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

(66) Para-fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide);

(67) Para-methylfentanyl (N-(4-methylphenyl)-N(1-phenethylpiperidin-4-yl)propionamide) (Other name: 4-methylfentanyl);

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

(69) Phenadoxone;

(70) Phenampromide;

(71) Phenomorphan;

(72) Phenoperidine;

(73) Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide) (Other name: benzoyl fentanyl);

(74) Piritramide;

(75) Proheptazine;

(76) Properidine;

(77) Propiram;

(78) Racemoramide;

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

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

(81) Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide) (Other names: 2-thiofuranyl fentanyl; thiophene fentanyl);

(82) Tilidine;

(83) Trimeperidine; [and]

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

2. By amending subsection (d) to read as follows:

"(d) Any material, compound, mixture, or preparation that contains any quantity of the following hallucinogenic substances, their salts, isomers, and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Alpha-ethyltryptamine (AET);

(2) 2,5-dimethoxy-4-ethylamphetamine (DOET);

(3) 2,5-dimethoxyamphetamine (2,5-DMA);

(4) 3,4-methylenedioxy amphetamine;

(5) 3,4-methylenedioxymethamphetamine (MDMA);

(6) N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-MDA);

(7) 3,4-methylenedioxy-N-ethylamphetamine (MDE);

(8) 5-methoxy-3,4-methylenedioxy-amphetamine;

(9) 4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);

(10) 4-Bromo-2,5-dimethoxyphenethylamine (Nexus);

(11) 3,4,5-trimethoxy amphetamine;

(12) Bufotenine;

(13) 4-methoxyamphetamine (PMA);

(14) Diethyltryptamine;

(15) Dimethyltryptamine;

(16) 4-methyl-2,5-dimethoxy-amphetamine;

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

(18) Ibogaine;

(19) Lysergic acid diethylamide;

(20) Marijuana;

(21) Parahexyl;

(22) Mescaline;

(23) Peyote;

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

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

(26) Psilocybin;

(27) Psilocyn;

(28) 1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);

(29) Ethylamine analog of phencyclidine (PCE);

(30) Pyrrolidine analog of phencyclidine (PCPy, PHP);

(31) Thiophene analog of phencyclidine (TPCP; TCP);

(32) Gamma-butyrolactone, including butyrolactone; butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone dihydro; dihydro-2(3H)furanone; tetrahydro-2-furanone; 1,2-butanolide; 1,4-butanolide; 4-butanolide; gamma-hydroxybutyric acid lactone; 3-hydroxybutyric acid lactone and 4-hydroxybutanoic acid lactone with Chemical Abstract Service number 96-48-0 when any such substance is intended for human ingestion;

(33) 1,4 butanediol, including butanediol; butane-1,4-diol; 1,4- butylenes glycol; butylene glycol; 1,4-dihydroxybutane; 1,4- tetramethylene glycol; tetramethylene glycol; tetramethylene 1,4- diol with Chemical Abstract Service number 110-63-4 when any such substance is intended for human ingestion;

(34) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7), its optical isomers, salts, and salts of isomers;

(35) N-benzylpiperazine (BZP; 1-benzylpiperazine) its optical isomers, salts, and salts of isomers;

(36) 1-(3-trifluoromethylphenyl)piperazine (TFMPP), its optical isomers, salts, and salts of isomers;

(37) Alpha-methyltryptamine (AMT), its isomers, salts, and salts of isomers;

(38) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT), its isomers, salts, and salts of isomers;

(39) Salvia divinorum;

(40) Salvinorin A;

(41) Divinorin A;

(42) 5-Methoxy-N,N-Dimethyltryptamine (5-MeO-DIPT) (some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT);

(43) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

(44) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);

(45) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

(46) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

(47) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2);

(48) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

(49) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);

(50) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);

(51) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);

(52) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5);

(53) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82); ~~and~~

(54) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);

(55) 2-(ethylamino)-2-(3-methoxylphenyl)cyclohexan-1-one(methoxetamine, MXE), 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;

(56) N-ethylhexedrone (Other names: aethylaminohexanophenone, 2-(ethylamino)-1-phenylhexan-s1-one);

(57) alpha-pyrrolidinohexanophenone (Other names: a-PHP, apyrrolidinohexanophenone, 1-phenyl-2-(pyrrolidine-1-yl)hexan-1-one);

(58) 4-methyl-alphaethylaminopentiophenone (Other names: 4-MEAP, 2-(ethylamino)-1-(4-methylphyl)pentan-1-one);

(59) 4'-methyl-alphapyrrolidinohexiophenone (Other names: MPHP, 4'-methyl-alphapyrrolidinohexanophenone, 1-(4-methylphenyl)-2-(pyrrolidine-1-yl)hexan1-one);

(60) Alpha-pyrrolidinoheptaphenone (Other names: PV8, 1-phenyl-2-(pyrrolidine-1-yl)heptan-1-one); and;

(61) 4'chloro-alphapyrrolidinovalerophenone (Other names: 4-chloro-a-PVP, 4'-chloro-apyrrolidinopentiophenone, 1-(4-chlorophenyl)-2-(pyrrolidine-1-yl)pentan-1-one) including their optical, positional, and geometric isomers, salts, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."

3. By amending subsection (e) to read as follows:

"(e) Depressants. Unless specifically excepted, the schedule shall include any material, compound, mixture, or preparation which contains any quantity of the substance:

(1) Mecloqualone;

(2) Methaqualone;

(3) Etizolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible); [or]

(4) Flualprazolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible)[.];

~~(5) N-ethylhexedrone (Other names: aethylaminohexanophenone, 2-(ethylamino)-1-phenylhexan-s1-one);~~

~~(6) alpha-pyrrolidinohexanophenone (Other names: a-PHP, apyrrolidinohexanophenone, 1-phenyl-2-(pyrrolidine-1-yl)hexan-1-one);~~

~~(7) 4-methyl-alphaethylaminopentiophenone (Other names: 4-MEAP, 2-(ethylamino)-1-(4-methylphyl)pentan-1-one);~~

~~(8) 4'-methyl-alphapyrrolidinohexiophenone (Other names: MPHP, 4'-methyl-alphapyrrolidinohexanophenone, 1-(4-methylphenyl)-2-(pyrrolidine-1-yl)hexan1-one);~~

~~(9) Alpha-pyrrolidinoheptaphenone (Other names: PV8, 1-phenyl-2-(pyrrolidine-1-yl)heptan-1-one); and~~

~~(10) 4'chloro-alphapyrrolidinovalerophenone (Other names: 4-chloro-a-PVP, 4'-chloro-apyrrolidinopentiophenone, 1-(4-chlorophenyl)-2-(pyrrolidine-1-yl)pentan-1-one) including their optical, positional, and geometric isomers, salts, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."~~

4. By amending subsection (f) to read as follows:

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

(1) Aminorex;

(2) Cathinone;

(3) 4,4'-dimethylaminorex (common name: 4,4'-DMAR);

(4) Fenethylline;

(5) Methcathinone;

(6) 4-methylaminorex;

(7) N-ethylamphetamine;

(8) N,N-dimethylamphetamine;

(9) Substituted cathinones, any compound, except bupropion or compounds listed under a different schedule, structurally derived from 2-

aminopropan-1-one by substitution at the 1-position with either phenyl, naphthyl, or thiophene ring systems, whether or not the compound is further modified 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; or

(C) 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.

Some other trade names: Mephedrone (2-methylamino-1-p-tolylpropan-1-one), also known as 4-methylmethcathinone (4-MMC), methylephedrone or MMCAT; Methylenedioxypyrovalerone (MDPV, MDPK); methylone or 3,4-methylenedioxymethcathinone; and 1-(benzo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one, monohydrochloride, also known as Ethylone, bk-MDEA hydrochloride, MDEC; 3,4-Methylenedioxy-N-ethylcathinone; bk-Methylenedioxyethylamphetamine, 4-methyl-N-ethylcathinone (4-MEC); 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP); alpha-pyrrolidinopentiophenone ([alpha]-PVP); 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-1-one (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone, bk-MBDP); 4-fluoro-N-methylcathinone (4-FMC, flephedrone); 3-fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-pyrrolidinobutiophenone ([alpha]-PBP) and their optical, positional, and geometric isomers, salts and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers is possible; and

(10) 1-(4-methoxyphenyl)-N-methylpropan-2-amine (Other names: para-methoxymethamphetamine; PMMA)[.];

~~(11)~~ Ethyl 2-(1-(5-fluoropentyl)-1Hindazole-3-carboxamido)-3,3-demethylbutanoate (Other name: 5FEDMB-PINACA);

~~(12)~~ Methyl 2-(1-(5-fluropentyl)-1Hindole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5FMDMB-PICA; 5F-MDMB-2201);

~~(13) N-(Adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-fluorobenzyl));~~

~~(14) 1-(5-Fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (Other names: 5F-CUMYLPINACA; SGT-25); and~~

~~(15) (1-(4-Fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144) including their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."~~

5. By amending subsection (g) to read as follows:

(g) Any of the following cannabinoids, their salts, isomers, and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Tetrahydrocannabinols; meaning tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. or synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following: Delta 1 cis or trans tetrahydrocannabinol, and their optical isomers; Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers; and 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, are covered);

(2) Naphthoylindoles; meaning any compound containing a 3-(1-naphthoyl)indole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 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 naphthyl ring to any extent;

(3) Naphthylmethylindoles; meaning any compound containing a 1H-indol-3-yl-(1-naphthyl) methane structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,

cycloalkylethyl, 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 naphthyl ring to any extent;

(4) Naphthoylpyrroles; meaning any compound containing a 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom of the pyrrole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the pyrrole ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(5) Naphthylmethylindenes; meaning any compound containing a naphthylideneindene structure with substitution at the 3-position of the indene ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the indene ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(6) Phenylacetylindoles; meaning any compound containing a 3-phenylacetylindole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the indole ring to any extent, whether or not substituted in the phenyl ring to any extent;

(7) Cyclohexylphenols; meaning any compound containing a 2-(3-hydroxycyclohexyl) phenol structure with substitution at the 5-position of the phenolic ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not substituted in the cyclohexyl ring to any extent;

(8) Benzoylindoles; meaning any compound containing a 3-(benzoyl) indole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 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;

(9) [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl) pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-naphthalenylmethanone (another trade name is WIN 55,212-2);

(10) (6a,10a)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol (Other trade names are: HU-210/HU-211);

(11) Tetramethylcyclopropanoylindoles; meaning any compound containing a 3-tetramethylcyclopropanoylindole structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the tetramethylcyclopropyl ring to any extent;

(12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: APINACA, AKB48);

(13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: PB-22; QUPIC);

(14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-carboxylate, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 5-fluoro-PB-22; 5F-PB-22);

(15) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-FUBINACA);

(16) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: ADB-PINACA);

(17) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-CHMINACA);

(18) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: AB-PINACA);

(19) [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl)methanone, and geometric isomers, salts, and salts of isomers (Other names: THJ-2201);

(20) Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate, and geometric isomers, salts, and salts of isomers (Other names: FUB-AMB, Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);

(21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, and geometric isomers, salts, and salts of isomers (Other names: 5-fluoro-AMB, 5-fluoro-AMP);

(22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: AKB48 N-(5-fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl analog, 5F-APINACA);

(23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and geometric isomers, salts, and salts of isomers (Other names: STS-135, 5F-APICA; 5-fluoro-APICA);

(24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate, and geometric isomers, salts, and salts of isomers (Other names: NM2201; CBL2201);

(25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide, and geometric isomers, salts, and salts of isomers (Other names: MAB-CHMINACA and ADB-CHMINACA);

(26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (Other names: 5F-ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical, positional, and geometric isomers, salts, and salts of isomers;

(27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA; CUMYL-4CN-BINACA);

(28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (Other name: 5F-AB-PINACA);

(29) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoate (Other names: MMB-CHMICA; AMB-CHMICA);

(30) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-CUMYL-P7AICA); ~~and~~

(31) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate (MDMB-4en-PINACA~~.~~;

(32) Ethyl 2-(1-(5-fluoropentyl)-1Hindazole-3-carboxamido)-3,3-demethylbutanoate (Other name: 5FEDMB-PINACA);

(33) Methyl 2-(1-(5-fluropentyl)-1Hindole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5FMDMB-PICA; 5F-MDMB-2201);

(34) N-(Adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-fluorobenzyl));

(35) 1-(5-Fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (Other names: 5F-CUMYLPINACA; SGT-25); and

(36) (1-(4-Fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144) including their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation."

SECTION 2. Section 329-20, Hawaii Revised Statutes, is amended by amending subsection (d) to read as follows:

1. By amending subsection (b) to read as follows:

"(b) Depressants. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, esters, ethers, and salts of isomers, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation, that has a degree of danger or probable danger associated with a depressant effect on the central nervous system:

(1) Alprazolam;

(2) Barbital;

(3) Brexanolone;

(4) Bromazepam;

(5) Butorphanol;

(6) Camazepam;

(7) Carisoprodol;

(8) Chloral betaine;

(9) Chloral hydrate;

(10) Chlordiazepoxide;

(11) Clobazam;

(12) Clonazepam;

(13) Clorazepate;

(14) Clotiazepam;

(15) Cloxazolam;

(16) Delorazepam;

(17) Diazepam;

(18) Dichloralphenazone (Midrin);

(19) Estazolam;

(20) Ethchlorvynol;

(21) Ethinamate;

(22) Ethyl loflazepate;

(23) Fludiazepam;

(24) Flunitrazepam;

(25) Flurazepam;

(26) Fospropofol (Lusedra);

(27) Halazepam;

(28) Haloxazolam;

(29) Ketazolam;

(30) Lemborexant ((1R,2S)-2-[(2,4-dimethylpyrimidin-5-yl)oxymethyl]-2-(3-fluorophenyl)-N-(5-fluoropyridin-2-yl)cyclopropane-1-carboxamide);

(31) Loprazolam;

(32) Lorazepam;

(33) Lormetazepam;

(34) Mebutamate;

(35) Medazepam;

(36) Meprobamate;

(37) Methohexital;

(38) Methylphenobarbital (mephorbarbital);

(39) Midazolam;

(40) Nimetazepam;

(41) Nitrazepam;

(42) Nordiazepam;

(43) Oxazepam;

(44) Oxazolam;

(45) Paraldehyde;

(46) Petrichloral;

(47) Phenobarbital;

(48) Pinazepam;

(49) Prazepam;

(50) Quazepam;

(51) Remimazolam;

(52) Suvorexant;

(53) Temazepam;

(54) Tetrazepam;

(55) Triazolam;

(56) Zaleplon;

(57) Zolpidem;

(58) Zopiclone (Lunesta); and

(59) Daridorexant and its salts."

2. By amending subsection (d) to read as follows:

"(d) Stimulants. Unless listed in another schedule, any material, compound, mixture, or preparation [which] that contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Cathine ((+)-norpseudoephedrine);

(2) Diethylpropion;

(3) Fencamfamin;

(4) Fenproporex;

(5) Lorcaserin;

(6) Mazindol;

(7) Mefenorex;

(8) Modafinil;

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

(10) Phentermine;

(11) Pipradrol;

(12) Serdexmethylphenidate;

(13) Sibutramine;

(14) Solriamfetol; [and]

(15) SPA (1-dimethylamino-1,2-dephenylethane, lefetamine)[.];

~~(16)~~ ~~Serdexmethylphenidate, including its salts, isomers, and salts of isomers; and~~

~~(17) Daridorexant and its salts.~~"

Thank you for the opportunity to testify in this measure.

HB-1097-HD-1
Submitted on: 2/22/2023 9:01:30 PM
Testimony for JHA on 2/23/2023 2:00:00 PM

Submitted By	Organization	Testifier Position	Testify
Johnnie-Mae L. Perry	Individual	Support	Written Testimony Only

Comments:

I, Johnnie-Mae L. Perry SUPPORT HB 1097 RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT.