THE SENATE
THIRTY-FIRST LEGISLATURE, 2022
STATE OF HAWAII

S.B. NO. 3141 S.D. 1 H.D. 2

A BILL FOR AN ACT

RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT.

BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF HAWAII:

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

1. By amending subsection (b) to read:

"(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);

[~~(3)~~] (6) Allylprodine;

[~~(4)~~] (7) Alphacetylmethadol (except levo-alphacetylmethadol, levomethadyl acetate, or LAAM);

[~~(5)~~] (8) Alphameprodine;

[~~(6)~~] (9) Alphamethadol;

[~~(7)~~] (10) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine);

[~~(8)~~] (11) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

[~~(9)~~] (12) Benzethidine;

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

[~~(10)~~] (14) Betacetylmethadol;

[~~(11)~~] (15) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide);

[~~(12)~~] (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);

[~~(13)~~] (18) Betameprodine;

[~~(14)~~] (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);

[~~(15)~~] (22) Betaprodine;

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

[~~(16)~~] (24) Clonitazene;

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

[~~(17)~~] (26) Dextromoramide;

[~~(18)~~] (27) Diampromide;

[~~(19)~~] (28) Diethylthiambutene;

[~~(20)~~] (29) Difenoxin;

[~~(21)~~] (30) Dimenoxadol;

[~~(22)~~] (31) Dimepheptanol;

[~~(23)~~] (32) Dimethylthiambutene;

[~~(24)~~] (33) Dioxaphetyl butyrate;

[~~(25)~~] (34) Dipipanone;

[~~(26)~~] (35) Ethylmethylthiambutene;

[~~(27)~~] (36) Etonitazene;

[~~(28)~~] (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);

[~~(29)~~] (42) Furethidine;

[~~(30)~~] (43) Hydroxypethidine;

[~~(31)~~] (44) Ketobemidone;

[~~(32)~~] (45) Levomoramide;

[~~(33)~~] (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);

[~~(34)~~] (49) [~~3-Methylfentanyl~~] 3-methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

[~~(35)~~] (50) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);

[~~(36)~~] (51) Morpheridine;

[~~(37)~~] (52) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);

[~~(38)~~] (53) Noracymethadol;

[~~(39)~~] (54) Norlevorphanol;

[~~(40)~~] (55) Normethadone;

[~~(41)~~] (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);

[~~(42)~~] (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);

[~~(43)~~] (68) PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine[~~,~~]);

[~~(44)~~] (69) Phenadoxone;

[~~(45)~~] (70) Phenampromide;

[~~(46)~~] (71) Phenomorphan;

[~~(47)~~] (72) Phenoperidine;

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

[~~(48)~~] (74) Piritramide;

[~~(49)~~] (75) Proheptazine;

[~~(50)~~] (76) Properidine;

[~~(51)~~] (77) Propiram;

[~~(52)~~] (78) Racemoramide;

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

[~~(53)~~] (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);

[~~(54)~~] (82) Tilidine;

[~~(55)~~] (83) Trimeperidine; and

[~~(56) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts, and salts of isomers;~~

~~(57) N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical isomers, salts, and salts of isomers;~~

~~(58) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide, (acetyl fentanyl), its optical, positional, and geometric isomers, salts, and salts of isomers;~~

~~(59) AH-7921 (3,4-dichloro-N-[(1-dimethylamino) cyclohexylmethyl]benzamide), its isomers, esters,~~

~~ethers, salts, and salts of isomers, esters, and ethers;~~

~~(60) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: Butyryl fentanyl);~~

~~(61) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, its isomers, esters, ethers, salts and salts of isomers, esters, and ethers (Other names: beta-hydroxythiofentanyl);~~

~~(62) N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: Furanyl fentanyl);~~

~~(63)~~] (84) U-47700 (3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide[~~, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: U-47700);~~

~~(64) 4-fluoroisobutyryl fentanyl or para-fluoroisobutyryl fentanyl [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide];~~

~~(65) Acryl fentanyl or acryloylfentanyl [N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide];~~

~~(66) Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide];~~

~~(67) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide;~~

~~(68) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);~~

~~(69) Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide) (Other name: 2-fluorofentanyl); and~~

~~(70) Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide~~]."

2. By amending subsections (f) and (g) to read:

"(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);

[~~(3)~~] (4) Fenethylline;

[~~(4)~~] (5) Methcathinone;

[~~(5) N-ethylamphetamine;~~]

(6) 4-methylaminorex;

(7) N-ethylamphetamine;

[~~(7)~~] (8) N,N-dimethylamphetamine; [~~and~~]

[~~(8)~~] (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).

(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; [~~and~~]

(27) [~~1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)indazole-3-carboxamide (CUMYL-4CN-BINACA),~~] 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its

optical, positional, and geometric isomers, salts, and salts of isomers[~~; also known as~~] (Other names: SGT-78[~~, 4-CN-CUMYL-BINACA~~]; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-CB-PINACA; CUMYL-CYBINACA; [~~4-cyano-CUMYL-BUTINACA.~~] 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)."

SECTION 2. Section 329-16, Hawaii Revised Statutes, is amended by amending subsection (c) to read as follows:

"(c) Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, whenever

the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

(1) Alfentanil;

(2) Alphaprodine;

(3) Anileridine;

(4) Bezitramide;

(5) Bulk Dextropropoxyphene (nondosage form);

(6) Carfentanil;

(7) Dihydrocodeine;

(8) Diphenoxylate;

(9) Fentanyl;

(10) Isomethadone;

(11) Levo-alphacetylmethadol (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-diphenyl-propane-carboxylic acid;

(18) Oliceridine, including the free base form, and its salts, to include the fumarate salt, by definition;

[~~(18)~~] (19) Pethidine (Meperidine);

[~~(19)~~] (20) Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine;

[~~(20)~~] (21) Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;

[~~(21)~~] (22) Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;

[~~(22)~~] (23) Phenazocine;

[~~(23)~~] (24) Piminodine;

[~~(24)~~] (25) Racemethorphan;

[~~(25)~~] (26) Racemorphan;

[~~(26)~~] (27) Remifentanil;

[~~(27)~~] (28) Sufentanil;

[~~(28)~~] (29) Tapentadol; and

[~~(29)~~] (30) Thiafentanil."

SECTION 3. Section 329-20, Hawaii Revised Statutes, is amended as follows:

1. By amending subsection (b) to read:

"(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;

[~~(3)~~] (4) Bromazepam;

[~~(4)~~] (5) Butorphanol;

[~~(5)~~] (6) Camazepam;

[~~(6)~~] (7) Carisoprodol;

[~~(7)~~] (8) Chloral betaine;

[~~(8)~~] (9) Chloral hydrate;

[~~(9)~~] (10) Chlordiazepoxide;

[~~(10)~~] (11) Clobazam;

[~~(11)~~] (12) Clonazepam;

[~~(12)~~] (13) Clorazepate;

[~~(13)~~] (14) Clotiazepam;

[~~(14)~~] (15) Cloxazolam;

[~~(15)~~] (16) Delorazepam;

[~~(16) Dichloralphenazone (Midrin);~~]

(17) Diazepam;

(18) Dichloralphenazone (Midrin);

[~~(18)~~] (19) Estazolam;

[~~(19)~~] (20) Ethchlorvynol;

[~~(20)~~] (21) Ethinamate;

[~~(21)~~] (22) Ethyl loflazepate;

[~~(22)~~] (23) Fludiazepam;

[~~(23)~~] (24) Flunitrazepam;

[~~(24)~~] (25) Flurazepam;

[~~(25)~~] (26) Fospropofol (Lusedra);

[~~(26)~~] (27) Halazepam;

[~~(27)~~] (28) Haloxazolam;

[~~(28)~~] (29) Ketazolam;

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

[~~(29)~~] (31) Loprazolam;

[~~(30)~~] (32) Lorazepam;

[~~(31)~~] (33) Lormetazepam;

[~~(32)~~] (34) Mebutamate;

[~~(33)~~] (35) Medazepam;

[~~(34)~~] (36) Meprobamate;

[~~(35)~~] (37) Methohexital;

[~~(36)~~] (38) Methylphenobarbital (mephorbarbital);

[~~(37)~~] (39) Midazolam;

[~~(38)~~] (40) Nimetazepam;

[~~(39)~~] (41) Nitrazepam;

[~~(40)~~] (42) Nordiazepam;

[~~(41)~~] (43) Oxazepam;

[~~(42)~~] (44) Oxazolam;

[~~(43)~~] (45) Paraldehyde;

[~~(44)~~] (46) Petrichloral;

[~~(45)~~] (47) Phenobarbital;

[~~(46)~~] (48) Pinazepam;

[~~(47)~~] (49) Prazepam;

[~~(48)~~] (50) Quazepam;

(51) Remimazolam;

[~~(49)~~] (52) Suvorexant;

[~~(50)~~] (53) Temazepam;

[~~(51)~~] (54) Tetrazepam;

[~~(52)~~] (55) Triazolam;

[~~(53)~~] (56) Zaleplon;

[~~(54)~~] (57) Zolpidem; and

[~~(55)~~] (58) Zopiclone (Lunesta)[~~; and~~

~~(56) Brexanolone~~]."

2. By amending subsection (d) to read:

"(d) Stimulants. 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 such 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;

[~~(5)~~] (6) Mazindol;

[~~(6)~~] (7) Mefenorex;

[~~(7)~~] (8) Modafinil;

[~~(8) Phentermine;~~]

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

(10) Phentermine;

[~~(10)~~] (11) Pipradrol;

(12) Serdexmethylphenidate;

[~~(11)~~] (13) Sibutramine;

(14) Solriamfetol; and

[~~(12)~~] (15) SPA (1-dimethylamino-1,2-diphenylethane, lefetamine) [~~;~~

~~(13) Lorcaserin; and~~

~~(14) Solriamfetol~~]."

SECTION 4. Section 329-22, Hawaii Revised Statutes, is amended by amending subsection (d) to read as follows:

"(d) Depressants. Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation that contains any quantity of the following substances having a depressant effect on the

central nervous system, including its salts, isomers, and salts of isomers:

(1) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (Other names: BRV; UCB-34714; Briviact) and its salts;

[~~(1)~~] (2) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide], (Vimpat);

(3) Lasmiditan (2,4,6-trifluoro-N-(6-(1-methylpiperidine-4-carbonyl)pyridine-2-yl-benzamide); and

[~~(2)~~] (4) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid][~~; and~~

~~(3) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (Other names: BRV; UCB-34714; Briviact) and its salts~~]."

SECTION 5. Statutory material to be repealed is bracketed and stricken. New statutory material is underscored.

SECTION 6. This Act shall take effect on January 1, 2050.

S.B. NO. 3141 S.D. 1 H.D. 2

Report Title:
Uniform Controlled Substances Act

Description:
Amends the schedules for controlled substances under the Uniform Controlled Substances Act for purposes of conforming with updates in federal law. Effective 1/1/2050. (HD2)

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