
A BILL FOR AN ACT

RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT.

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

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

3 1. By amending subsection (d) to read:

4 "(d) ~~[Any]~~ Hallucinogenic substances. Unless specifically
5 excepted or unless listed in another schedule, any material,
6 compound, mixture, or preparation that contains any quantity of
7 the following hallucinogenic substances, including their salts,
8 isomers, and salts of isomers, ~~[unless specifically excepted,]~~
9 whenever the existence of these salts, isomers, and salts of
10 isomers is possible within the specific chemical designation:

- 11 (1) Alpha-ethyltryptamine (AET);
12 (2) 2,5-dimethoxy-4-ethylamphetamine (DOET);
13 (3) 2,5-dimethoxyamphetamine (2,5-DMA);
14 (4) 3,4-methylenedioxy amphetamine;
15 (5) 3,4-methylenedioxymethamphetamine (MDMA);
16 (6) N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-
17 MDA);



- 1 (7) 3,4-methylenedioxy-N-ethylamphetamine (MDE);
- 2 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;
- 3 (9) 4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);
- 4 (10) 4-Bromo-2,5-dimethoxyphenethylamine (Nexus);
- 5 (11) 3,4,5-trimethoxy amphetamine;
- 6 (12) Bufotenine;
- 7 (13) 4-methoxyamphetamine (PMA);
- 8 (14) Diethyltryptamine;
- 9 (15) Dimethyltryptamine;
- 10 (16) 4-methyl-2,5-dimethoxy-amphetamine;
- 11 (17) Gamma hydroxybutyrate (GHB) (some other names include
- 12 gamma hydroxybutyric acid; 4-hydroxybutyrate; 4-
- 13 hydroxybutanoic acid; sodium oxybate; sodium
- 14 oxybutyrate);
- 15 (18) Ibogaine;
- 16 (19) Lysergic acid diethylamide;
- 17 (20) Marijuana;
- 18 (21) Parahexyl;
- 19 (22) Mescaline;
- 20 (23) Peyote;
- 21 (24) N-ethyl-3-piperidyl benzilate;



- 1 (25) N-methyl-3-piperidyl benzilate;
- 2 (26) Psilocybin;
- 3 (27) Psilocyn;
- 4 (28) 1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);
- 5 (29) Ethylamine analog of phencyclidine (PCE);
- 6 (30) Pyrrolidine analog of phencyclidine (PCPy, PHP);
- 7 (31) Thiophene analog of phencyclidine (TCP; TCP);
- 8 (32) Gamma-butyrolactone, including butyrolactone;
- 9 butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone
- 10 dihydro; dihydro-2(3H)furanone; tetrahydro-2-furanone;
- 11 1,2-butanolide; 1,4-butanolide; 4-butanolide; gamma-
- 12 hydroxybutyric acid lactone; 3-hydroxybutyric acid
- 13 lactone and 4-hydroxybutanoic acid lactone with
- 14 Chemical Abstract Service number 96-48-0 when any such
- 15 substance is intended for human ingestion;
- 16 (33) 1,4 butanediol, including butanediol; butane-1,4-diol;
- 17 1,4- butylenes glycol; butylene glycol; 1,4-
- 18 dihydroxybutane; 1,4- tetramethylene glycol;
- 19 tetramethylene glycol; tetramethylene 1,4- diol with
- 20 Chemical Abstract Service number 110-63-4 when any
- 21 such substance is intended for human ingestion;



- 1 (34) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7),
2 its optical isomers, salts, and salts of isomers;
- 3 (35) N-benzylpiperazine (BZP; 1-benzylpiperazine), its
4 optical isomers, salts, and salts of isomers;
- 5 (36) 1-(3-trifluoromethylphenyl)piperazine (TFMPP), its
6 optical isomers, salts, and salts of isomers;
- 7 (37) Alpha-methyltryptamine (AMT) [~~its isomers, salts, and~~
8 ~~salts of isomers~~];
- 9 (38) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT) [~~its~~
10 ~~isomers, salts, and salts of isomers~~];
- 11 (39) Salvia divinorum;
- 12 (40) Salvinorin A;
- 13 (41) Divinorin A;
- 14 (42) 5-Methoxy-N,N-Dimethyltryptamine (5-MeO-DIPT) (some
15 trade or other names: 5-methoxy-3-[2-
16 (dimethylamino)ethyl]indole; 5-MeO-DMT);
- 17 (43) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);
- 18 (44) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);
- 19 (45) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);
- 20 (46) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);



- 1 (47) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-
2 2);
- 3 (48) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine
4 (2C-T-4);
- 5 (49) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);
- 6 (50) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);
- 7 (51) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);
- 8 (52) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-
9 methoxybenzyl)ethanamine, its optical, positional, and
10 geometric isomers, salts, and salts of isomers (Other
11 names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5);
- 12 (53) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-
13 methoxybenzyl)ethanamine, its optical, positional, and
14 geometric isomers, salts, and salts of isomers (Other
15 names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82); [and]
- 16 (54) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-
17 methoxybenzyl)ethanamine, its optical, positional, and
18 geometric isomers, salts, and salts of isomers (Other
19 names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36) [-];
- 20 (55) N-ethylhexedrone, its optical, positional, and
21 geometric isomers, salts, and salts of isomers (Other



- 1 names: [alpha]-ethylaminohexanophenone; 2-
2 (ethylamino)-1-phenylhexan-1-one);
- 3 (56) Alpha-pyrrolidinohexanophenone, its optical,
4 positional, and geometric isomers, salts, and salts of
5 isomers (Other names: [alpha]-PHP; [alpha]-
6 pyrrolidinohexanophenone; 1-phenyl-2-(pyrrolidin-1-
7 yl)hexan-1-one);
- 8 (57) 4-methyl-alpha-ethylaminopentiophenone, its optical,
9 positional, and geometric isomers, salts, and salts of
10 isomers (Other names: 4-MEAP; 2-(ethylamino)-1-(4-
11 methylphenyl)pentan-1-one);
- 12 (58) 4'-methyl-alpha-pyrrolidinohexiophenone, its optical,
13 positional, and geometric isomers, salts, and salts of
14 isomers (Other names: MPHP; 4'-methyl-alpha-
15 pyrrolidinohexanophenone; 1-(4-methylphenyl)-2-
16 (pyrrolidin-1-yl)hexan-1-one);
- 17 (59) Alpha-pyrrolidinoheptaphenone, its optical,
18 positional, and geometric isomers, salts, and salts of
19 isomers (Other names: PV8; 1-phenyl-2-(pyrrolidin-1-
20 yl)heptan-1-one);



- 1 (60) 4'-chloro-alpha-pyrrolidinovalerophenone, its optical,
2 positional, and geometric isomers, salts, and salts of
3 isomers (Other names: 4-chloro-[alpha]-PVP; 4'-
4 chloro-[alpha]-pyrrolidinopentiophenone; 1-(4-
5 chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one); and
6 (61) 2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one
7 (methoxetamine, MXE)."

8 2. By amending subsection (g) to read:

9 "(g) [Any] Cannabinoids. Unless specifically excepted or
10 unless listed in another schedule, any of the following
11 cannabinoids, including their salts, isomers, and salts of
12 isomers, [~~unless specifically excepted,~~] whenever the existence
13 of these salts, isomers, and salts of isomers is possible within
14 the specific chemical designation:

- 15 (1) Tetrahydrocannabinols; meaning tetrahydrocannabinols
16 naturally contained in a plant of the genus Cannabis
17 (cannabis plant), as well as synthetic equivalents of
18 the substances contained in the plant, or in the
19 resinous extractives of Cannabis, sp. or synthetic
20 substances, derivatives, and their isomers with
21 similar chemical structure and pharmacological



1 activity to those substances contained in the plant,
2 such as the following: Delta 1 cis or trans
3 tetrahydrocannabinol, and their optical isomers; Delta
4 6 cis or trans tetrahydrocannabinol, and their optical
5 isomers; and Delta 3,4 cis or trans-
6 tetrahydrocannabinol, and its optical isomers (since
7 nomenclature of these substances is not
8 internationally standardized, compounds of these
9 structures, regardless of numerical designation of
10 atomic positions, are covered);

11 (2) Naphthoylindoles; meaning any compound containing a 3-
12 (1-naphthoyl)indole structure with substitution at the
13 nitrogen atom of the indole ring by a alkyl,
14 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
15 1-(N-methyl-2-piperidiny)methyl or 2-(4-
16 morpholinyl)ethyl group, whether or not further
17 substituted in the indole ring to any extent and
18 whether or not substituted in the naphthyl ring to any
19 extent;

20 (3) Naphthylmethylinindoles; meaning any compound containing
21 a 1H-indol-3-yl-(1-naphthyl) methane structure with



1 substitution at the nitrogen atom of the indole ring
2 by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
3 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
4 2-(4-morpholinyl) ethyl group whether or not further
5 substituted in the indole ring to any extent and
6 whether or not substituted in the naphthyl ring to any
7 extent;

8 (4) Naphthoylpyrroles; meaning any compound containing a
9 3-(1-naphthoyl)pyrrole structure with substitution at
10 the nitrogen atom of the pyrrole ring by a alkyl,
11 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)
13 ethyl group whether or not further substituted in the
14 pyrrole ring to any extent, whether or not substituted
15 in the naphthyl ring to any extent;

16 (5) Naphthylmethylindenes; meaning any compound containing
17 a naphthylideneindene structure with substitution at
18 the 3-position of the indene ring by a alkyl,
19 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
20 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)
21 ethyl group whether or not further substituted in the



1 indene ring to any extent, whether or not substituted
2 in the naphthyl ring to any extent;

3 (6) Phenylacetylindoles; meaning any compound containing a
4 3-phenylacetylindole structure with substitution at
5 the nitrogen atom of the indole ring by a alkyl,
6 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
7 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)
8 ethyl group whether or not further substituted in the
9 indole ring to any extent, whether or not substituted
10 in the phenyl ring to any extent;

11 (7) Cyclohexylphenols; meaning any compound containing a
12 2-(3-hydroxycyclohexyl) phenol structure with
13 substitution at the 5-position of the phenolic ring by
14 a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
15 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
16 2-(4-morpholinyl) ethyl group whether or not
17 substituted in the cyclohexyl ring to any extent;

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



1 1-(N-methyl-2-piperidinyl) methyl, or 2-(4-
2 morpholinyl) ethyl group whether or not further
3 substituted in the indole ring to any extent and
4 whether or not substituted in the phenyl ring to any
5 extent;

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

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

14 (11) Tetramethylcyclopropanoylindoles; meaning any compound
15 containing a 3-tetramethylcyclopropanoylindole
16 structure with substitution at the nitrogen atom of
17 the indole ring by an alkyl, haloalkyl, cyanoalkyl,
18 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
19 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,
20 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
21 morpholinyl)methyl, or tetrahydropyranylmethyl group,



- 1 whether or not further substituted in the indole ring
2 to any extent and whether or not substituted in the
3 tetramethylcyclopropyl ring to any extent;
- 4 (12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
5 its optical, positional, and geometric isomers, salts,
6 and salts of isomers (Other names: APINACA, AKB48);
- 7 (13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
8 optical, positional, and geometric isomers, salts, and
9 salts of isomers (Other names: PB-22; QUPIC);
- 10 (14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-
11 carboxylate, its optical, positional, and geometric
12 isomers, salts, and salts of isomers (Other names: 5-
13 fluoro-PB-22; 5F-PB-22);
- 14 (15) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-
15 fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
16 positional, and geometric isomers, salts, and salts of
17 isomers (Other names: AB-FUBINACA);
- 18 (16) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
19 indazole-3-carboxamide, its optical, positional, and
20 geometric isomers, salts, and salts of isomers (Other
21 names: ADB-PINACA);



- 1 (17) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
2 (cyclohexylmethyl)-1H-indazole-3-carboxamide, its
3 optical, positional, and geometric isomers, salts, and
4 salts of isomers (Other names: AB-CHMINACA);
- 5 (18) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
6 indazole-3-carboxamide, and geometric isomers, salts,
7 and salts of isomers (Other names: AB-PINACA);
- 8 (19) [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-
9 yl)methanone, and geometric isomers, salts, and salts
10 of isomers (Other names: THJ-2201);
- 11 (20) Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-
12 valinate, and geometric isomers, salts, and salts of
13 isomers (Other names: FUB-AMB, Methyl 2-(1-(4-
14 fluorobenzyl)-1H-indazole-3-carboxamido)-3-
15 methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);
- 16 (21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
17 carboxamido)-3-methylbutanoate, and geometric isomers,
18 salts, and salts of isomers (Other names: 5-fluoro-
19 AMB, 5-fluoro-AMP);
- 20 (22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-
21 indazole-3-carboxamide, and geometric isomers, salts,



- 1 and salts of isomers (Other names: AKB48 N-(5-
2 fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
3 analog, 5F-APINACA);
- 4 (23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
5 geometric isomers, salts, and salts of isomers (Other
6 names: STS-135, 5F-APICA; 5-fluoro-APICA);
- 7 (24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
8 carboxylate, and geometric isomers, salts, and salts
9 of isomers (Other names: NM2201; CBL2201);
- 10 (25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
11 (cyclohexylmethyl)-1H-indazole-3-carboxamide, and
12 geometric isomers, salts, and salts of isomers (Other
13 names: MAB-CHMINACA and ADB-CHMINACA);
- 14 (26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
15 carboxamido]-3,3-dimethylbutanoate (Other names: 5F-
16 ADB, [~~5-fluoro-ADB~~], 5-fluoro-ADB, and 5F-MDMB-
17 PINACA), its optical, positional, and geometric
18 isomers, salts, and salts of isomers;
- 19 (27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-
20 3-carboxamide, its optical, positional, and geometric
21 isomers, salts, and salts of isomers (Other names:



- 1 SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-
- 2 CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA;
- 3 CUMYL-4CN-BINACA) ;
- 4 (28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
- 5 fluoropentyl)-1H-indazole-3-carboxamide (Other name:
- 6 5F-AB-PINACA) ;
- 7 (29) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
- 8 carboxamido)-3-methylbutanoate (Other names: MMB-
- 9 CHMICA; AMB-CHMICA) ;
- 10 (30) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
- 11 pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-
- 12 CUMYL-P7AICA) ; [and]
- 13 (31) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
- 14 3-carboxamido)butanoate (MDMB-4en-PINACA) [-] ;
- 15 (32) Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
- 16 carboxamido)-3,3-dimethylbutanoate (Other name: 5F-
- 17 EDMB-PINACA) ;
- 18 (33) Methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-
- 19 3,3-dimethylbutanoate (Other names: 5F-MDMB-PICA; 5F-
- 20 MDMB-2201) ;



- 1 (34) N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-
2 carboxamide (Other names: FUB-AKB48; FUB-APINACA;
3 AKB48 N-(4-FLUOROBENZYL));
- 4 (35) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
5 indazole-3-carboxamide (Other names: 5F-CUMYL-PINACA;
6 SGT-25); and
- 7 (36) (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-
8 tetramethylcyclopropyl)methanone (Other name: FUB-
9 144)."

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

12 "(b) Depressants. [~~Any~~] Unless specifically excepted or
13 unless listed in another schedule, any material, compound,
14 mixture, or preparation [which] that contains any quantity of
15 the following substances, including its salts, isomers, esters,
16 ethers, and salts of isomers, whenever the existence of these
17 isomers, esters, ethers, and salts is possible within the
18 specific chemical designation, that has a degree of danger or
19 probable danger associated with a depressant effect on the
20 central nervous system:

21 (1) Alprazolam;



- 1 (2) Barbital;
- 2 (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) Daridorexant;
- 16 [~~+16~~] (17) Delorazepam;
- 17 [~~+17~~] (18) Diazepam;
- 18 [~~+18~~] (19) Dichloralphenazone (Midrin);
- 19 [~~+19~~] (20) Estazolam;
- 20 [~~+20~~] (21) Ethchlorvynol;
- 21 [~~+21~~] (22) Ethinamate;



- 1 [~~(22)~~] (23) Ethyl loflazepate;
- 2 [~~(23)~~] (24) Fludiazepam;
- 3 [~~(24)~~] (25) Flunitrazepam;
- 4 [~~(25)~~] (26) Flurazepam;
- 5 [~~(26)~~] (27) Fospropofol (Lusedra);
- 6 [~~(27)~~] (28) Halazepam;
- 7 [~~(28)~~] (29) Haloxazolam;
- 8 [~~(29)~~] (30) Ketazolam;
- 9 [~~(30)~~] (31) Lemborexant ((1R,2S)-2-[(2,4-dimethylpyrimidin-
10 5-yl)oxymethyl]-2-(3-fluorophenyl)-N-(5-fluoropyridin-
11 2-yl)cyclopropane-1-carboxamide);
- 12 [~~(31)~~] (32) Loprazolam;
- 13 [~~(32)~~] (33) Lorazepam;
- 14 [~~(33)~~] (34) Lormetazepam;
- 15 [~~(34)~~] (35) Mebutamate;
- 16 [~~(35)~~] (36) Medazepam;
- 17 [~~(36)~~] (37) Meprobamate;
- 18 [~~(37)~~] (38) Methohexital;
- 19 [~~(38)~~] (39) Methylphenobarbital (mephobarbital);
- 20 [~~(39)~~] (40) Midazolam;
- 21 [~~(40)~~] (41) Nimetazepam;



1 [~~(41)~~] (42) Nitrazepam;
2 [~~(42)~~] (43) Nordiazepam;
3 [~~(43)~~] (44) Oxazepam;
4 [~~(44)~~] (45) Oxazolam;
5 [~~(45)~~] (46) Paraldehyde;
6 [~~(46)~~] (47) Petrichloral;
7 [~~(47)~~] (48) Phenobarbital;
8 [~~(48)~~] (49) Pinazepam;
9 [~~(49)~~] (50) Prazepam;
10 [~~(50)~~] (51) Quazepam;
11 [~~(51)~~] (52) Remimazolam;
12 [~~(52)~~] (53) Suvorexant;
13 [~~(53)~~] (54) Temazepam;
14 [~~(54)~~] (55) Tetrazepam;
15 [~~(55)~~] (56) Triazolam;
16 [~~(56)~~] (57) Zaleplon;
17 [~~(57)~~] (58) Zolpidem; and
18 [~~(58)~~] (59) Zopiclone (Lunesta)."

19 SECTION 3. Section 329-22, Hawaii Revised Statutes, is
20 amended by amending subsection (d) to read as follows:



1 "(d) Depressants. Unless specifically exempted or
2 excluded or unless listed in another schedule, any material,
3 compound, mixture, or preparation that contains any quantity of
4 the following substances having a depressant effect on the
5 central nervous system, including its salts, isomers, and salts
6 of isomers:

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

10 (2) Ganaxolone (3[alpha]-hydroxy-3[beta]-methyl-5[alpha]-
11 pregnan-20-one);

12 ~~[(2)]~~ (3) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-
13 propionamide], (Vimpat);

14 ~~[(3)]~~ (4) Lasmiditan (2,4,6-trifluoro-N-(6-(1-
15 methylpiperidine-4-carbonyl)pyridine-2-yl)-benzamide);

16 and

17 ~~[(4)]~~ (5) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic
18 acid]."

19 SECTION 4. Statutory material to be repealed is bracketed
20 and stricken. New statutory material is underscored.



1 SECTION 5. This Act shall take effect upon its approval.



Report Title:

Uniform Controlled Substances Act

Description:

Updates the Uniform Controlled Substances Act by adding substances scheduled under federal law. (SD1)

The summary description of legislation appearing on this page is for informational purposes only and is not legislation or evidence of legislative intent.

