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: 1. By amending subsection (d) to read: 3 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 $(\dot{2})$ 2,5-dimethoxy-4-ethylamphetamine (DOET); 13 2,5-dimethoxyamphetamine (2,5-DMA); (3) 14 3,4-methylenedioxy amphetamine; (4)15 (5) 3,4-methylenedioxymethamphetamine (MDMA); 16 N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-(6) 17 MDA);

```
3,4-methylenedioxy-N-ethylamphetamine (MDE);
 1
          (7)
               5-methoxy-3,4-methylenedioxy-amphetamine;
 2
         (8)
               4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);
 3
         (9)
 4
         (10)
               4-Bromo-2,5-dimethoxyphenethylamine (Nexus);
               3,4,5-trimethoxy amphetamine;
 5
         (11)
 6
         (12)
              Bufotenine;
 7
               4-methoxyamphetamine (PMA);
         (13)
 8
         (14)
              Diethyltryptamine;
 9
              Dimethyltryptamine;
         (15)
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
              Ibogaine;
         (18)
              Lysergic acid diethylamide;
16
         (19)
17
        (20)
              Marijuana;
              Parahexyl;
18
         (21)
19
         (22) Mescaline;
20
        (23) Peyote;
21
              N-ethyl-3-piperidyl benzilate;
         (24)
```

```
(25)
              N-methyl-3-piperidyl benzilate;
 1
 2
              Psilocybin;
        (26)
 3
        (27)
              Psilocyn;
              1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);
 4
        (28)
              Ethylamine analog of phencyclidine (PCE);
 5
        (29)
              Pyrrolidine analog of phencyclidine (PCPy, PHP);
 6
        (30)
 7
              Thiophene analog of phencyclidine (TPCP; TCP);
        (31)
 8
        (32)
              Gamma-butyrolactone, including butyrolactone;
              butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone
 9
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;
```

H.B. NO. 1097 H.D. 2

```
1
        (34)
              2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7),
              its optical isomers, salts, and salts of isomers;
 2
 3
        (35)
              N-benzylpiperazine (BZP; 1-benzylpiperazine), its
              optical isomers, salts, and salts of isomers;
 4
 5
              1-(3-trifluoromethylphenyl)piperazine (TFMPP), its
        (36)
 6
              optical isomers, salts, and salts of isomers;
 7
              Alpha-methyltryptamine (AMT) [, its isomers, salts, and
        (37)
 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
              2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);
        (43)
18
        (44)
              2-(2,5-Dimethoxy-4-methylphenyl) ethanamine (2C-D);
              2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C);
19
        (45)
20
        (46)
              2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);
```

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

1		names: [alpha]-ethylaminohexanophenone; 2-
2		<pre>(ethylamino) -1-phenylhexan-1-one);</pre>
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		<pre>methylphenyl)pentan-1-one);</pre>
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 li	sted in another schedule, any of the following
11	cannabino	ids, <u>including</u> 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 speci	fic 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-piperidinyl)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)	Naphthylmethylindoles; meaning any compound containing
21		a 1H-indol-3-yl-(1-naphthyl) methane structure with

1		substitution at the nitrogen atom of the indoie 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

H.B. NO. 1097 H.D. 2

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-(N-methyl-2-piperidinyl) methyl, or 2-(4-
 1
              morpholinyl) ethyl group whether or not further
 2
              substituted in the indole ring to any extent and
3
              whether or not substituted in the phenyl ring to any
 4
 5
               extent;
6
         (9)
              [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
              pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-
7
               naphthalenylmethanone (another trade name is WIN
8
               55,212-2);
9
10
               (6a, 10a) - 9 - (hydroxymethyl) - 6, 6 - dimethyl - 3 - (2 - a)
        (10)
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
               structure with substitution at the nitrogen atom of
16
17
              the indole ring by an alkyl, haloalkyl, cyanoalkyl,
18
              alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
              methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl,
19
20
              1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
              morpholinyl) methyl, or tetrahydropyranylmethyl group,
21
```

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

```
N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
 1
         (17)
               (cyclohexylmethyl) -1H-indazole-3-carboxamide, its
 2
               optical, positional, and geometric isomers, salts, and
 3
               salts of isomers (Other names: AB-CHMINACA);
 4
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
 5
        (18)
               indazole-3-carboxamide, and geometric isomers, salts,
 6
               and salts of isomers (Other names: AB-PINACA);
 7
 8
               [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-
        (19)
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-
               fluorobenzyl)-1H-indazole-3-carboxamido)-3-
14
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,
```

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

H.B. NO. 1097 H.D. 2

```
SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-
1
              CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA;
2
3
              CUMYL-4CN-BINACA);
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
4
        (28)
              fluoropentyl) -1H-indazole-3-carboxamide (Other name:
5
 6
              5F-AB-PINACA);
7
              Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
        (29)
              carboxamido) - 3-methylbutanoate (Other names: MMB-
8
              CHMICA; AMB-CHMICA);
9
10
              1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
        (30)
11
              pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-
12
              CUMYL-P7AICA); [and]
              Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
13
        (31)
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		<u>144).</u> "	
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.li	sted 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;	

```
Barbital;
 1
          (2)
                Brexanolone;
 2
          (3)
 3
          (4)
                Bromazepam;
                Butorphanol;
 4
          (5)
 5
          (6)
               Camazepam;
 6
                Carisoprodol;
          (7)
               Chloral betaine;
 7
          (8)
 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
        [\frac{(17)}{(18)}] (18) Diazepam;
18
        [<del>(18)</del>] (19) Dichloralphenazone (Midrin);
19
        [<del>(19)</del>] (20) Estazolam;
20
        [\frac{(20)}{}] (21) Ethchlorvynol;
21
        \left[\frac{(21)}{(21)}\right] (22) Ethinamate;
```

```
[\frac{(22)}{(23)}] (23) Ethyl loflazepate;
 1
 2
          \left[\frac{(23)}{(24)}\right] (24) Fludiazepam;
 3
          [<del>(24)</del>] (25) Flunitrazepam;
 4
          \left[\frac{(25)}{(25)}\right] (26) Flurazepam;
          [\frac{(26)}{(27)}] (27) Fospropofol (Lusedra);
 5
 6
          \left[\frac{(27)}{}\right] (28) Halazepam;
 7
          \left[\frac{(28)}{(29)}\right] (29) Haloxazolam;
 8
          [\frac{(29)}{(30)}] (30) Ketazolam;
 9
          [\frac{(30)}{(31)}] (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
          [\frac{(32)}{(33)}] (33) Lorazepam;
14
          [(33)] (34) Lormetazepam;
15
          \left[\frac{(34)}{(35)}\right] (35) Mebutamate;
16
          [-(35)] (36) Medazepam;
17
          [\frac{(36)}{(37)}] (37) Meprobamate;
18
          [\frac{(37)}{}] (38) Methohexital;
19
          [<del>(38)</del>] (39) Methylphenobarbital (mephorbarbital);
20
          \left[\frac{(39)}{(40)}\right] (40) Midazolam;
21
          [(40)] (41) Nimetazepam;
```

```
1
          \left[\frac{(41)}{(41)}\right] (42) Nitrazepam;
          \left[\frac{(42)}{(43)}\right] (43) Nordiazepam;
 2
          [-(43)-] (44) Oxazepam;
 3
          [-(44)] (45) Oxazolam;
 4
          [<del>(45)</del>] (46) Paraldehyde;
 5
 6
          [<del>(46)</del>] (47) Petrichloral;
          [(47)) (48) Phenobarbital;
 7
          \left[\frac{(48)}{(49)}\right] (49) Pinazepam;
 9
          [-(49)] (50) Prazepam;
10
          [-(50)] (51) Quazepam;
11
          [\frac{(51)}{(52)}] Remimazolam;
12
          \left[\frac{(52)}{(53)}\right] (53) Suvorexant;
13
          \left[\frac{(53)}{(54)}\right] (54) Temazepam;
14
         \left[\frac{(54)}{(55)}\right] (55) Tetrazepam;
15
          \left[\frac{(55)}{(56)}\right] (56) Triazolam;
          [\frac{(56)}{(57)}] (57) Zaleplon;
16
          [\frac{(57)}{(58)}] Zolpidem; and
17
          [<del>(58)</del>] (59) Zopiclone (Lunesta)."
18
19
             SECTION 3. Section 329-22, Hawaii Revised Statutes, is
      amended by amending subsection (d) to read as follows:
20
```

```
"(d) Depressants. Unless specifically exempted or
 1
     excluded or unless listed in another schedule, any material,
 2
    compound, mixture, or preparation that contains any quantity of
3
     the following substances having a depressant effect on the
4
     central nervous system, including its salts, isomers, and salts
 5
6
    of isomers:
          (1) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-
7
               yl]butanamide) (Other names: BRV; UCB-34714;
8
9
               Briviact) and its salts;
          (2) Ganaxolone (3[alpha]-hydroxy-3[beta]-methyl-5[alpha]-
10
11
               pregnan-20-one);
12
         \left[\frac{(2)}{(2)}\right] (3) Lacosamide \left[\frac{(R)}{2}\right] -2-acetoamido-N-benzyl-3-methoxy-
13
               propionamide], (Vimpat);
         \left[\frac{3}{3}\right] (4) Lasmiditan (2,4,6-trifluoro-N-(6-(1-
14
15
               methylpiperidine-4-carbonyl)pyridine-2-yl-benzamide);
               and
16
         \left[\frac{4}{4}\right] (5) Pregabalin \left[\frac{5}{3}\right] (aminomethyl) -5-methylhexanoic
17
18
               acid]."
          SECTION 4. Statutory material to be repealed is bracketed
19
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. (HD2)

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