S.B. NO. 2882 S.D. 2 H.D. 1

# A BILL FOR AN ACT

RELATING TO THE UNIFORM CONTROLLED SUBSTANCES ACT.

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

| 1 | SECTIO     | N 1.  | Section | 329-14, | Hawaii | Revised | Statutes, | is |
|---|------------|-------|---------|---------|--------|---------|-----------|----|
| 2 | amended as | follo | ws:     |         |        |         |           |    |

| 3 | 1. | Ву | amending | subsection | (d) | to | read: |
|---|----|----|----------|------------|-----|----|-------|
|---|----|----|----------|------------|-----|----|-------|

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

- 10 (1) Alpha-ethyltryptamine (AET);
- 11 (2) 2,5-dimethoxy-4-ethylamphetamine (DOET);
- 12 (3) 2,5-dimethoxyamphetamine (2,5-DMA);
- 13 (4) 3,4-methylenedioxy amphetamine;
- 14 (5) 3,4-methylenedioxymethamphetamine (MDMA);
- 15 (6) N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-16 MDA);
- 17 (7) 3,4-methylenedioxy-N-ethylamphetamine (MDE);
- 18 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;

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| 1  | (9)  | 4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);  |
|----|------|---|
| 2  | (10) | 4-Bromo-2,5-dimethoxyphenethylamine (Nexus);          |
| 3  | (11) | 3,4,5-trimethoxy amphetamine;                         |
| 4  | (12) | Bufotenine;   |
| 5  | (13) | 4-methoxyamphetamine (PMA);                           |
| 6  | (14) | Diethyltryptamine;                                    |
| 7  | (15) | Dimethyltryptamine;                                   |
| 8  | (16) | 4-methyl-2,5-dimethoxy-amphetamine;                   |
| 9  | (17) | Gamma hydroxybutyrate (GHB) (some other names include |
| 10 |      | gamma hydroxybutyric acid; 4-hydroxybutyrate; 4-      |
| 11 |      | hydroxybutanoic acid; sodium oxybate; sodium          |
| 12 |      | oxybutyrate);   |
| 13 | (18) | Ibogaine;   |
| 14 | (19) | Lysergic acid diethylamide;                           |
| 15 | (20) | Marijuana;  |
| 16 | (21) | Parahexyl;  |
| 17 | (22) | Mescaline;  |
| 18 | (23) | Peyote;   |
| 19 | (24) | N-ethyl-3-piperidyl benzilate;                        |
| 20 | (25) | N-methyl-3-piperidyl benzilate;                       |
| 21 | (26) | Psilocybin;   |
| 22 | (27) | Psilocyn;   |

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| 1  | (28) | 1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);       |
|----|------|--|
| 2  | (29) | Ethylamine analog of phencyclidine (PCE);              |
| 3  | (30) | Pyrrolidine analog of phencyclidine (PCPy, PHP);       |
| 4  | (31) | Thiophene analog of phencyclidine (TPCP; TCP);         |
| 5  | (32) | Gamma-butyrolactone, including butyrolactone;          |
| 6  |      | butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone   |
| 7  |      | dihydro; dihydro-2(3H)furanone; tetrahydro-2-furanone; |
| 8  |      | 1,2-butanolide; 1,4-butanolide; 4-butanolide; gamma-   |
| 9  |      | hydroxybutyric acid lactone; 3-hydroxybutyric acid     |
| 10 |      | lactone and 4-hydroxybutanoic acid lactone with        |
| 11 |      | Chemical Abstract Service number 96-48-0 when any such |
| 12 |      | substance is intended for human ingestion;             |
| 13 | (33) | 1,4 butanediol, including butanediol; butane-1,4-diol; |
| 14 |      | 1,4- butylenes glycol; butylene glycol; 1,4-           |
| 15 |      | dihydroxybutane; 1,4- tetramethylene glycol;           |
| 16 |      | tetramethylene glycol; tetramethylene 1,4- diol with   |
| 17 |      | Chemical Abstract Service number 110-63-4 when any     |
| 18 |      | such substance is intended for human ingestion;        |
| 19 | (34) | 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7), |
| 20 |      | its optical isomers, salts, and salts of isomers;      |
| 21 | (35) | N-benzylpiperazine (BZP; 1-benzylpiperazine) its       |
| 22 |      | optical isomers, salts, and salts of isomers;          |
|    |      |  |



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| 1           | (36) | 1-(3-trifluoromethylphenyl)piperazine (TFMPP), its     |
|-------------|------|--|
| 2           |      | optical isomers, salts, and salts of isomers;          |
| 3           | (37) | Alpha-methyltryptamine (AMT), its isomers, salts, and  |
| 4           |      | salts of isomers;                                      |
| 5           | (38) | 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT), its  |
| 6           |      | isomers, salts, and salts of isomers;                  |
| 7           | (39) | Salvia divinorum;                                      |
| 8           | (40) | Salvinorin A;  |
| 9           | (41) | Divinorin A; [ <del>and</del> ]                        |
| 10          | (42) | 5-Methoxy-N,N-Dimethyltryptamine (5-MeO-DIPT) (some    |
| 11          |      | trade or other names: 5-methoxy-3-[2-                  |
| 12          |      | (dimethylamino)ethyl]indole; 5-MeO-DMT)[+];            |
| 13          | (43) | 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);      |
| 14          | (44) | 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D);     |
| 15          | (45) | 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);     |
| <b>16</b> · | (46) | 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);       |
| 17          | (47) | 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T- |
| 18          |      | <u>2);</u>   |
| 19          | (48) | 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine    |
| 20          |      | <u>(2C-T-4);</u>                                       |
| 21          | (49) | 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);              |
| 22          | (50) | 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N);     |
|             |      |  |



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| 1  | (51)       | 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P);  |
|----|------------|---|
| 2  | (52)       | 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-                    |
| 3  |            | methoxybenzyl)ethanamine, its optical, positional, and  |
| 4  |            | geometric isomers, salts, and salts of isomers (Other   |
| 5  |            | <pre>names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5);</pre> |
| 6  | (53)       | 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-                  |
| 7  |            | methoxybenzyl)ethanamine, its optical, positional, and  |
| 8  |            | geometric isomers, salts, and salts of isomers (Other   |
| 9  |            | names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82); and       |
| 10 | (54)       | 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-                   |
| 11 |            | methoxybenzyl)ethanamine, its optical, positional, and  |
| 12 |            | geometric isomers, salts, and salts of isomers (Other   |
| 13 |            | names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36)."          |
| 14 | 2. 1       | By amending subsection (g) to read:                     |
| 15 | " (g)      | Any of the following cannabinoids, their salts,         |
| 16 | isomers, a | and salts of isomers, unless specifically excepted,     |
| 17 | whenever   | the existence of these salts, isomers, and salts of     |
| 18 | isomers is | s possible within the specific chemical designation:    |
| 19 | (1)        | Tetrahydrocannabinols; meaning tetrahydrocannabinols    |
| 20 |            | naturally contained in a plant of the genus Cannabis    |
| 21 |            | (cannabis plant), as well as synthetic equivalents of   |
| 22 |            | the substances contained in the plant, or in the        |
|    | an0.000    |   |



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1 resinous extractives of Cannabis, sp. or synthetic substances, derivatives, and their isomers with 2 similar chemical structure and pharmacological 3 4 activity to those substances contained in the plant, 5 such as the following: Delta 1 cis or trans 6 tetrahydrocannabinol, and their optical isomers; Delta 7 6 cis or trans tetrahydrocannabinol, and their optical 8 isomers; and Delta 3,4 cis or trans-9 tetrahydrocannabinol, and its optical isomers (since 10 nomenclature of these substances is not 11 internationally standardized, compounds of these 12 structures, regardless of numerical designation of 13 atomic positions, are covered); 14 (2) Naphthoylindoles; meaning any compound containing a 15 3-(1-naphthoyl) indole structure with substitution at 16 the nitrogen atom of the indole ring by a alkyl, 17 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) 18 19 ethyl group, whether or not further substituted in the 20 indole ring to any extent and whether or not 21 substituted in the naphthyl ring to any extent;



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

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

18 (5) Naphthylmethylindenes; meaning any compound containing
19 a naphthylideneindene structure with substitution at
20 the 3-position of the indene ring by a alkyl,
21 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,

22 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)



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| 1  |           | ethyl group whether or not further substituted in the  |
|----|-----------|--|
| 2  |           | indene ring to any extent, whether or not substituted  |
| 3  |           | in the naphthyl ring to any extent;                    |
| 4  | (6)       | Phenylacetylindoles; meaning any compound containing a |
| 5  |           | 3-phenylacetylindole structure with substitution at    |
| 6  |           | the nitrogen atom of the indole ring by a alkyl,       |
| 7  |           | haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, |
| 8  |           | 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) |
| 9  |           | ethyl group whether or not further substituted in the  |
| 10 |           | indole ring to any extent, whether or not substituted  |
| 11 |           | in the phenyl ring to any extent;                      |
| 12 | (7)       | Cyclohexylphenols; meaning any compound containing a   |
| 13 |           | 2-(3-hydroxycyclohexyl) phenol structure with          |
| 14 |           | substitution at the 5-position of the phenolic ring by |
| 15 |           | a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,         |
| 16 |           | cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or  |
| 17 |           | 2-(4-morpholinyl) ethyl group whether or not           |
| 18 |           | substituted in the cyclohexyl ring to any extent;      |
| 19 | (8)       | Benzoylindoles; meaning any compound containing a      |
| 20 |           | 3-(benzoyl) indole structure with substitution at the  |
| 21 |           | nitrogen atom of the indole ring by a alkyl,           |
| 22 |           | haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, |
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| 1  |      | 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)          |
|----|------|---|
| 2  |      | ethyl group whether or not further substituted in the           |
| 3  |      | indole ring to any extent and whether or not                    |
| 4  |      | substituted in the phenyl ring to any extent;                   |
| 5  | (9)  | 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)                    |
| 6  |      | pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-                       |
| 7  |      | napthalenylmethanone (another trade name is WIN                 |
| 8  |      | 55,212-2);  |
| 9  | (10) | (6a,10a)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-                  |
| 10 |      | methyloctan-2-yl)-6a,7,10,10a-                                  |
| 11 |      | tetrahydrobenzo[c]chromen-1-ol (other trade names are:          |
| 12 |      | HU-210 and HU-211); [and]                                       |
| 13 | (11) | Tetramethylcyclopropanoylindoles; meaning any compound          |
| 14 |      | containing a 3-tetramethylcyclopropanoylindole                  |
| 15 |      | structure with substitution at the nitrogen atom of             |
| 16 |      | the indole ring by an alkyl, haloalkyl, cyanoalkyl,             |
| 17 |      | alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-               |
| 18 |      | <pre>methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,</pre> |
| 19 |      | 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-               |
| 20 |      | morpholinyl)methyl, or tetrahydropyranylmethyl group,           |
| 21 |      | whether or not further substituted in the indole ring           |



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| 1  |      | to any extent and whether or not substituted in the    |
|----|------|--|
| 2  |      | tetramethylcyclopropyl ring to any $extent[-]$ ;       |
| 3  | (12) | N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,    |
| 4  |      | its optical, positional, and geometric isomers, salts, |
| 5  |      | and salts of isomers (Other names: APINACA, AKB48);    |
| 6  | (13) | Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its    |
| 7  |      | optical, positional, and geometric isomers, salts, and |
| 8  |      | salts of isomers (Other names: PB-22; QUPIC);          |
| 9  | (14) | <u>Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-</u>    |
| 10 |      | carboxylate, its optical, positional, and geometric    |
| 11 |      | isomers, salts, and salts of isomers (Other names: 5-  |
| 12 |      | fluoro-PB-22; 5F-PB-22);                               |
| 13 | (15) | N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-             |
| 14 |      | fluorobenzyl)-1H-indazole-3-carboxamide, its optical,  |
| 15 |      | positional, and geometric isomers, salts, and salts of |
| 16 |      | isomers (Other names: AB-FUBINACA); and                |
| 17 | (16) | N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-  |
| 18 |      | indazole-3-carboxamide, its optical, positional, and   |
| 19 |      | geometric isomers, salts, and salts of isomers (Other  |
| 20 |      | names: ADB-PINACA)."                                   |
| 21 | SECT | ION 2. Section 329-18, Hawaii Revised Statutes, is     |

22 amended by amending subsection (c) to read as follows:



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"(c) Depressants. Unless listed in another schedule, any
 material, compound, mixture, or preparation containing any
 quantity of the following substances having a depressant effect
 on the central nervous system:

5 Any compound, mixture, or preparation containing (1) 6 amobarbital, secobarbital, pentobarbital, or any salt 7 thereof and one or more other active medicinal 8 ingredients which are not listed in any schedule; 9 (2) Any suppository dosage form containing amobarbital, 10 secobarbital, pentobarbital, or any salt of any of 11 these drugs and approved by the Food and Drug 12 Administration for marketing only as a suppository; 13 (3) Any substance that contains any quantity of a 14 derivative of barbituric acid or any salt thereof, 15 including the substance butalbital; 16 (4)Chlorhexadol; 17 (5) Embutramide (Tributame);

- 18 (6) Ketamine, its salts, isomers, and salts of isomers,
- 19 also known as (+ or -)-2-(2-chlorophenyl)-2-
- 20 (methylamino)-cyclohexanone;
- 21 (7) Lysergic acid;
- 22 (8) Lysergic acid amide;



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| 1  | (9)   | Methyprylon;  |
|----|---|---|
| 2  | (10)  | Sulfondiethylmethane;                                 |
| 3  | (11)  | Sulfonethylmethane;                                   |
| 4  | (12)  | Sulfonmethane;  |
| 5  | (13)  | Tiletamine/Zolazepam (Telazol, 2-(ethylamino)-2-(-    |
| 6  |   | thienyl)-cyclohexanone, flupyrazapon) or any salts    |
| 7  |   | thereof; [and]  |
| 8  | (14)  | Gamma hydroxybutyric acid and its salts, isomers, and |
| 9  |   | salts of isomers that are contained in a drug product |
| 10 |   | for which an application has been approved under      |
| 11 |   | section 505 of the federal Food, Drug, and Cosmetic   |
| 12 |   | Act[+];   |
| 13 | (15)  | Perampanel [2-(2-oxo-1-phenyl-5-pyridin-2-yl-1,2-     |
| 14 |   | dihydropyridin-3-yl)benzonitrile], including its      |
| 15 |   | salts, isomers, and salts of isomers; and             |
| 16 | (16)  | Alfaxalone 5[alpha]-pregnan-3[alpha]-ol-11,20-dione,  |
| 17 |   | including its salts, isomers, and salts of isomers."  |
| 18 | SECT  | ION 3. Section 329-20, Hawaii Revised Statutes, is    |
| 19 | amended by                                  | y amending subsection (d) to read as follows:         |
| 20 | "(d)  | Stimulants. Unless listed in another schedule, any    |
| 21 | material,                                   | compound, mixture, or preparation which contains any  |
| 22 | quantity o                                  | of the following substances having a stimulant effect |
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| 1  | on the ce | entral nervous system, including its salts, isomers, and |
|----|-----------|--|
| 2  | salts of  | such isomers whenever the existence of such salts,       |
| 3  | isomers,  | and salts of isomers is possible within the specific     |
| 4  | chemical  | designation:   |
| 5  | (1)       | Cathine ((+)-norpseudoephedrine);                        |
| 6  | (2)       | Diethylpropion;  |
| 7  | (3)       | Fencamfamin;   |
| 8  | (4)       | Fenproporex;   |
| 9  | (5)       | Mazindol;  |
| 10 | (6)       | Mefenorex;   |
| 11 | (7)       | Modafinil;   |
| 12 | (8)       | Phentermine;   |
| 13 | (9)       | Pemoline (including organometallic complexes and         |
| 14 |           | chelates thereof);                                       |
| 15 | (10)      | Pipradrol;   |
| 16 | (11)      | Sibutramine; [and]                                       |
| 17 | (12)      | SPA (1-dimethylamino-1,2-diphenylethane,                 |
| 18 |           | lefetamine)[-]; and                                      |
| 19 | (13)      | Lorcaserin."   |
| 20 | SECT      | ION 4. Statutory material to be repealed is bracketed    |
| 21 | and stric | ken. New statutory material is underscored.              |
| 22 | SECT      | ION 5. This Act shall take effect on July 1, 2112.       |

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#### Report Title:

Uniform Controlled Substances Act; Federal Conformity

#### Description:

SB2882 HD1 HMS 2014-2537

Updates chapter 329, Hawaii Revised Statutes (HRS), to make it consistent with amendments to federal law on controlled substances. Amends sections 329-14, 329-18, and 329-20, HRS, to add new controlled substances federally scheduled as required under section 329-11, HRS. Effective July 1, 2112. (SB2882 HD1)

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