

JAN 23 2015

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# A BILL FOR AN ACT

RELATING TO THE SCHEDULING OF MARIJUANA.

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

1           SECTION 1. The legislature finds that the medical use of  
2 marijuana has existed in the State since 2000 when Hawaii became  
3 the first state to accept the medical use of marijuana by the  
4 legislative process. However, despite fourteen years of medical  
5 use by thousands of patients in the State, marijuana still  
6 remains in schedule I, while coca and poppy, plants that contain  
7 substances far more dangerous than any substance found in  
8 marijuana, are listed in schedule II.

9           The legislature further finds that Hawaii's administrative  
10 process for making recommendations to the legislature on the  
11 scheduling of controlled substances is devoid of any scientific  
12 or medical input from the department of health. Therefore,  
13 statutory changes are needed in order to include the department  
14 of health in this process.

15           The purpose of this Act is to remove marijuana and its  
16 intrinsic cannabinoids from schedule I and require the  
17 department of public safety, in consultation with the department



1 of health, to issue a recommendation for the rescheduling of  
2 marijuana.

3 SECTION 2. Section 329-11, Hawaii Revised Statutes, is  
4 amended by amending subsections (a) and (b) to read as follows:

5 "(a) Annually, upon the convening of each regular session  
6 of the state legislature, the department of public safety shall  
7 report to the legislature additions, deletions, or revisions in  
8 the schedules of substances enumerated in sections 329-14,  
9 329-16, 329-18, 329-20, and 329-22, and any other  
10 recommendations that it deems necessary. Three months prior to  
11 the convening of each regular session, the department of public  
12 safety shall post public notice, at the state capitol and in the  
13 office of the lieutenant governor for public inspection, of the  
14 department's recommendations to the legislature concerning any  
15 additions, deletions, or revisions in these schedules; provided  
16 that the posting shall not be required if official notice has  
17 been received that the substance has been added, deleted, or  
18 rescheduled as a controlled substance under federal law. In  
19 making a determination regarding a substance, the department of  
20 public safety, in consultation with the department of health,



1 shall assess the degree of danger or probable danger of the  
2 substance by considering the following:

3 (1) The actual or probable abuse of the substance  
4 including:

5 (A) Its history and current pattern of abuse;

6 (B) The scope, duration, and significance of abuse;  
7 and

8 (C) A judgment of the degree of actual or probable  
9 detriment that may result from the abuse of the  
10 substance;

11 (2) The biomedical hazard of the substance including:

12 (A) Its pharmacology: the effects and modifiers of  
13 effects of the substance;

14 (B) Its toxicology: the acute and chronic toxicity,  
15 interaction with other substances whether  
16 controlled or not, and liability to psychic or  
17 physiological dependence;

18 (C) Risk to public health and particular  
19 susceptibility of segments of the population; and



(D) Existence of therapeutic alternatives for substances that are or may be used for medical purposes;

(3) A judgment of the probable physical and social impact of widespread abuse of the substance;

(4) Whether the substance is an immediate precursor of a substance already controlled under this part; and

(5) The current state of scientific knowledge regarding the substance.

(b) After considering a scientific and medical evaluation from the department of health regarding the factors enumerated in subsection (a) [7] and conducting formal rulemaking proceedings pursuant to chapter 91, the department of public safety shall make a recommendation to the legislature, specifying to what schedule the substance should be added, deleted, or rescheduled if it finds that the substance has a degree of danger or probable danger. The evaluation and recommendations of the department of health shall be made in writing and submitted to the department of public safety no later than October 1, 2015. The recommendations of the department of health shall be binding on the department of



1 public safety as to such scientific and medical matters. If the  
2 department of health recommends that a drug or other substance  
3 not be controlled, the department of public safety shall not  
4 recommend control of the drug or other substance.

5 Recommendations of the department of health shall become  
6 effective immediately. The department of public safety may make  
7 its recommendation to the legislature prior to the submission of  
8 its annual report, in which case the department of public safety  
9 shall publish and give notice to the public of the  
10 recommendation."

11 SECTION 3. Section 329-14, Hawaii Revised Statutes, is  
12 amended as follows:

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

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

20 (1) Alpha-ethyltryptamine (AET);

21 (2) 2,5-dimethoxy-4-ethylamphetamine (DOET);



- 1 (3) 2,5-dimethoxyamphetamine (2,5-DMA);
- 2 (4) 3,4-methylenedioxy amphetamine;
- 3 (5) 3,4-methylenedioxymethamphetamine (MDMA);
- 4 (6) N-hydroxy-3,4-methylenedioxyamphetamine (N-hydroxy-
- 5 MDA);
- 6 (7) 3,4-methylenedioxy-N-ethylamphetamine (MDE);
- 7 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;
- 8 (9) 4-bromo-2,5-dimethoxy-amphetamine (4-bromo-2,5-DMA);
- 9 (10) 4-Bromo-2,5-dimethoxyphenethylamine (Nexus);
- 10 (11) 3,4,5-trimethoxy amphetamine;
- 11 (12) Bufotenine;
- 12 (13) 4-methoxyamphetamine (PMA);
- 13 (14) Diethyltryptamine;
- 14 (15) Dimethyltryptamine;
- 15 (16) 4-methyl-2,5-dimethoxy-amphetamine;
- 16 (17) Gamma hydroxybutyrate (GHB) (some other names include
- 17 gamma hydroxybutyric acid; 4-hydroxybutyrate;
- 18 4-hydroxybutanoic acid; sodium oxybate; sodium
- 19 oxybutyrate);
- 20 (18) Ibogaine;
- 21 (19) Lysergic acid diethylamide;



1        [~~(20)~~] ~~Marijuana~~;  
2        ~~(21)~~] (20)    Parahexyl;  
3        [~~(22)~~] (21)    Mescaline;  
4        [~~(23)~~] (22)    Peyote;  
5        [~~(24)~~] (23)    N-ethyl-3-piperidyl benzilate;  
6        [~~(25)~~] (24)    N-methyl-3-piperidyl benzilate;  
7        [~~(26)~~] (25)    Psilocybin;  
8        [~~(27)~~] (26)    Psilocyn;  
9        [~~(28)~~] (27)    1-[1-(2-Thienyl) cyclohexyl] Pyrrolidine (TCPy);  
10       [~~(29)~~] (28)    Ethylamine analog of phencyclidine (PCE);  
11       [~~(30)~~] (29)    Pyrrolidine analog of phencyclidine (PCPy, PHP);  
12       [~~(31)~~] (30)    Thiophene analog of phencyclidine (TPCP; TCP);  
13       [~~(32)~~] (31)    Gamma-butyrolactone, including butyrolactone;  
14           butyrolactone gamma; 4-butyrolactone; 2(3H)-furanone  
15           dihydro; dihydro-2(3H)furanone; tetrahydro-2-furanone;  
16           1,2-butanolide; 1,4-butanolide; 4-butanolide; gamma-  
17           hydroxybutyric acid lactone; 3-hydroxybutyric acid  
18           lactone and 4-hydroxybutanoic acid lactone with  
19           Chemical Abstract Service number 96-48-0 when any such  
20           substance is intended for human ingestion;



- 1       ~~[(+33+)]~~ (32) 1,4 butanediol, including butanediol; butane-  
2               1,4-diol; 1,4- butylenes glycol; butylene glycol;  
3               1,4-dihydroxybutane; 1,4- tetramethylene glycol;  
4               tetramethylene glycol; tetramethylene 1,4- diol with  
5               Chemical Abstract Service number 110-63-4 when any  
6               such substance is intended for human ingestion;
- 7       ~~[(+34+)]~~ (33) 2,5-dimethoxy-4-(n)-propylthiophenethylamine  
8               (2C-T-7), its optical isomers, salts, and salts of  
9               isomers;
- 10       ~~[(+35+)]~~ (34) N-benzylpiperazine (BZP; 1-benzylpiperazine) its  
11               optical isomers, salts, and salts of isomers;
- 12       ~~[(+36+)]~~ (35) 1-(3-trifluoromethylphenyl)piperazine (TFMPP),  
13               its optical isomers, salts, and salts of isomers;
- 14       ~~[(+37+)]~~ (36) Alpha-methyltryptamine (AMT), its isomers,  
15               salts, and salts of isomers;
- 16       ~~[(+38+)]~~ (37) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-  
17               DIPT), its isomers, salts, and salts of isomers;
- 18       ~~[(+39+)]~~ (38) Salvia divinorum;
- 19       ~~[(+40+)]~~ (39) Salvinorin A;
- 20       ~~[(+41+)]~~ (40) Divinorin A;





- 1        [~~(42)~~] (41)    5-Methoxy-N,N-Dimethyltryptamine (5-MeO-DIPT)  
2                    (some trade or other names: 5-methoxy-3-[2-  
3                    (dimethylamino)ethyl]indole; 5-MeO-DMT);  
4        [~~(43)~~] (42)    2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine  
5                    (2C-E);  
6        [~~(44)~~] (43)    2-(2,5-Dimethoxy-4-methylphenyl)ethanamine  
7                    (2C-D);  
8        [~~(45)~~] (44)    2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine  
9                    (2C-C);  
10       [~~(46)~~] (45)    2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);  
11       [~~(47)~~] (46)    2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine  
12                    (2C-T-2);  
13       [~~(48)~~] (47)    2-[4-(Isopropylthio)-2,5-  
14                    dimethoxyphenyl]ethanamine (2C-T-4);  
15       [~~(49)~~] (48)    2-(2,5-Dimethoxyphenyl)ethanamine (2C-H);  
16       [~~(50)~~] (49)    2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine  
17                    (2C-N);  
18       [~~(51)~~] (50)    2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine  
19                    (2C-P);  
20       [~~(52)~~] (51)    2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-  
21                    methoxybenzyl)ethanamine, its optical, positional, and



1 geometric isomers, salts, and salts of isomers (Other  
2 names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5);

3 ~~[(+53)]~~ (52) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-  
4 methoxybenzyl)ethanamine, its optical, positional, and  
5 geometric isomers, salts, and salts of isomers (Other  
6 names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82); and

7 ~~[(+54)]~~ (53) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-  
8 methoxybenzyl)ethanamine, its optical, positional, and  
9 geometric isomers, salts, and salts of isomers (Other  
10 names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36)."

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

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

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



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

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

20 ~~[(3)]~~ (2) Naphthylmethylinindoles; meaning any compound  
21 containing a 1H-indol-3-yl-(1-naphthyl) methane



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

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

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



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

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

11 ~~[+7+]~~ (6) Cyclohexylphenols; meaning any compound  
12 containing a 2-(3-hydroxycyclohexyl) phenol structure  
13 with substitution at the 5-position of the phenolic  
14 ring by 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+]~~ (7) Benzoylindoles; meaning any compound containing a  
19 3-(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-piperidiny1) 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+]~~ (8) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)  
6 pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-  
7 naphthalenylmethanone (another trade name is WIN  
8 55,212-2);

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

13 ~~[+11+]~~ (10) Tetramethylcyclopropanoylindoles; meaning any  
14 compound containing a 3-  
15 tetramethylcyclopropanoylindole structure with  
16 substitution at the nitrogen atom of the indole ring  
17 by an alkyl, haloalkyl, cyanoalkyl, alkenyl,  
18 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-  
19 piperidiny1)methyl, 2-(4-morpholinyl)ethyl, 1-(N-  
20 methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
21 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)]~~ (11) 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)]~~ (12) 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)]~~ (13) 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)]~~ (14) 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); and

~~[(16)]~~ (15) 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)."

SECTION 4. The department of public safety, in  
consultation with the department of health, shall make its  
recommendations for the rescheduling of marijuana and  
tetrahydrocannabinols no later than October 1, 2015. These  
recommendations shall become effective immediately.

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

SECTION 6. This Act shall take effect on July 1, 2015.

INTRODUCED BY:

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# S.B. NO. 189

**Report Title:**

Marijuana; Rescheduling; Department of Health; Department of Public Safety

**Description:**

Removes marijuana and its intrinsic cannabinoids from schedule I and requires the department of public safety, in consultation with the department of health, to issue a recommendation for the rescheduling of marijuana.

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

