



EXECUTIVE CHAMBERS  
HONOLULU

DAVID Y. IGE  
GOVERNOR

July 10, 2018

**GOV. MSG. NO. 1299**

The Honorable Ronald D. Kouchi,  
President  
and Members of the Senate  
Twenty-Ninth State Legislature  
State Capitol, Room 409  
Honolulu, Hawai'i 96813

The Honorable Scott K. Saiki,  
Speaker and Members of the  
House of Representatives  
Twenty-Ninth State Legislature  
State Capitol, Room 431  
Honolulu, Hawai'i 96813

Dear President Kouchi, Speaker Saiki, and Members of the Legislature:

This is to inform you that on July 10, 2018, the following bill was signed into law:

HB2385 HD2 SD1

RELATING TO THE UNIFORM CONTROLLED  
SUBSTANCES ACT  
**ACT 190 (18)**

Sincerely,

DAVID Y. IGE  
Governor, State of Hawai'i

Approved by the Governor  
JUL 10 2018  
on \_\_\_\_\_

ORIGINAL

ACT 190

HOUSE OF REPRESENTATIVES  
TWENTY-NINTH LEGISLATURE, 2018  
STATE OF HAWAII

H.B. NO. 2385  
H.D. 2  
S.D. 1

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# 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 by amending subsections (f) and (g) to read as follows:

3 "(f) Stimulants. Unless specifically excepted or unless  
4 listed in another schedule, any material, compound, mixture, or  
5 preparation which contains any quantity of the following  
6 substances having a stimulant effect on the central nervous  
7 system, including its salts, isomers, and salts of isomers:

- 8 (1) Aminorex;  
9 (2) Cathinone;  
10 (3) Fenethylamine;  
11 (4) Methcathinone;  
12 (5) N-ethylamphetamine;  
13 (6) 4-methylaminorex;  
14 (7) N,N-dimethylamphetamine; and  
15 (8) Substituted cathinones, any compound, except bupropion  
16 or compounds listed under a different schedule,  
17 structurally derived from 2-aminopropan-1-ol by



1 substitution at the 1-position with either phenyl,  
2 naphthyl, or thiophene ring systems, whether or not  
3 the compound is further modified in any of the  
4 following ways:

- 5 (A) By substitution in the ring system to any extent  
6 with alkyl, alkylenedioxy, alkoxy, haloalkyl,  
7 hydroxyl, or halide substituents, whether or not  
8 further substituted in the ring system by one or  
9 more other univalent substituents;
- 10 (B) By substitution at the 3-position with an acyclic  
11 alkyl substituent; or
- 12 (C) By substitution at the 2-amino nitrogen atom with  
13 alkyl, dialkyl, benzyl, or methoxybenzyl groups,  
14 or by inclusion of the 2-amino nitrogen atom in a  
15 cyclic structure.

16 Some other trade names: Mephedrone (2-methylamino-1-  
17 p-tolylpropan-1-one), also known as 4-  
18 methylmethcathinone (4-MMC), methylephedrone or MMCAT;  
19 Methylenedioxypropylone (MDPV, MDPK); methylone or  
20 3,4-methylenedioxymethcathinone; and 1-  
21 (benzo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one,



1 monohydrochloride, also known as Ethylone, bk-MDEA  
2 hydrochloride, MDEC; 3,4-Methylenedioxy-N-  
3 ethylcathinone; bk-Methylenedioxyethylamphetamine [±],  
4 4-methyl-N-ethylcathinone (4-MEC); 4-methyl-alpha-  
5 pyrrolidinopropiophenone (4-MePPP); alpha-  
6 pyrrolidinopentiophenone ([alpha]-PVP); 1-(1,3-  
7 benzodioxol-5-yl)-2-(methylamino)butan-1-one  
8 (butylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-  
9 1-one (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-  
10 (methylamino)pentan-1-one (pentylone, bk-MBDP); 4-  
11 fluoro-N-methylcathinone (4-FMC, flephedrone); 3-  
12 fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-  
13 2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-  
14 pyrrolidinobutiophenone ([alpha]-PBP) and their  
15 optical, positional, and geometric isomers, salts and  
16 salts of isomers, whenever the existence of such  
17 salts, isomers, and salts of isomers is possible.

18 (g) Any of the following cannabinoids, their salts,  
19 isomers, and salts of isomers, unless specifically excepted,  
20 whenever the existence of these salts, isomers, and salts of  
21 isomers is possible within the specific chemical designation:



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



- 1 pyrrole ring to any extent, whether or not substituted  
2 in the naphthyl ring to any extent;
- 3 (5) Naphthylmethylindenes; meaning any compound containing  
4 a naphthylideneindene structure with substitution at  
5 the 3-position of the indene 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 indene ring to any extent, whether or not substituted  
10 in the naphthyl ring to any extent;
- 11 (6) Phenylacetylindoles; meaning any compound containing a  
12 3-phenylacetylindole structure with substitution at  
13 the nitrogen atom of the indole ring by a alkyl,  
14 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,  
15 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)  
16 ethyl group whether or not further substituted in the  
17 indole ring to any extent, whether or not substituted  
18 in the phenyl ring to any extent;
- 19 (7) Cyclohexylphenols; meaning any compound containing a  
20 2-(3-hydroxycyclohexyl) phenol structure with  
21 substitution at the 5-position of the phenolic ring by



- 1 a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,  
2 cycloalkylethyl, 1-(N-methyl-2-piperidiny) methyl or  
3 2-(4-morpholinyl) ethyl group whether or not  
4 substituted in the cyclohexyl ring to any extent;
- 5 (8) Benzoylindoles; meaning any compound containing a 3-  
6 (benzoyl) indole structure with substitution at the  
7 nitrogen atom of the indole ring by a alkyl,  
8 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,  
9 1-(N-methyl-2-piperidiny) methyl, or 2-(4-  
10 morpholinyl) ethyl group whether or not further  
11 substituted in the indole ring to any extent and  
12 whether or not substituted in the phenyl ring to any  
13 extent;
- 14 (9) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)  
15 pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-  
16 naphthalenylmethanone (another trade name is WIN  
17 55,212-2);
- 18 (10) (6a,10a)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-  
19 methyloctan-2-yl)-6a,7,10,10a-  
20 tetrahydrobenzo[c]chromen-1-ol (Other trade names are:  
21 HU-210/HU-211);





- 1 (11) Tetramethylcyclopropanoylindoles; meaning any compound  
2 containing a 3-tetramethylcyclopropanoylindole  
3 structure with substitution at the nitrogen atom of  
4 the indole ring by an alkyl, haloalkyl, cyanoalkyl,  
5 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-  
6 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,  
7 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
8 morpholinyl)methyl, or tetrahydropyranylmethyl group,  
9 whether or not further substituted in the indole ring  
10 to any extent and whether or not substituted in the  
11 tetramethylcyclopropyl ring to any extent;
- 12 (12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,  
13 its optical, positional, and geometric isomers, salts,  
14 and salts of isomers (Other names: APINACA, AKB48);
- 15 (13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its  
16 optical, positional, and geometric isomers, salts, and  
17 salts of isomers (Other names: PB-22; QUPIC);
- 18 (14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-  
19 carboxylate, its optical, positional, and geometric  
20 isomers, salts, and salts of isomers (Other names: 5-  
21 fluoro-PB-22; 5F-PB-22);



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



- 1 (21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-  
2 carboxamido)-3-methylbutanoate, and geometric isomers,  
3 salts, and salts of isomers (Other names: 5-fluoro-  
4 AMB, 5-fluoro-AMP);
- 5 (22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-  
6 indazole-3-carboxamide, and geometric isomers, salts,  
7 and salts of isomers (Other names: AKB48 N-(5-  
8 fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl  
9 analog, 5F-APINACA);
- 10 (23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and  
11 geometric isomers, salts, and salts of isomers (Other  
12 names: STS-135, 5F-APICA; 5-fluoro-APICA);
- 13 (24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-  
14 carboxylate, and geometric isomers, salts, and salts  
15 of isomers (Other names: NM2201);
- 16 (25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-  
17 (cyclohexylmethyl)-1H-indazole-3-carboxamide, and  
18 geometric isomers, salts, and salts of isomers (Other  
19 names: MAB-CHMINACA and ADB-CHMINACA); [and]
- 20 (26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-  
21 carboxamido]-3,3-dimethylbutanoate (Other names: 5F-



1 ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical,  
2 positional, and geometric isomers, salts, and salts of  
3 isomers [±]; and

4 (27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)indazole-3-  
5 carboxamide (CUMYL-4CN-BINACA), its optical,  
6 positional, and geometric isomers, salts, and salts of  
7 isomers; also known as SGT-78, 4-CN-CUMYL-BINACA;  
8 CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano CUMYL-  
9 BUTINACA."

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

12 "(g) Hallucinogenic substances, unless listed in another  
13 schedule, shall include:

14 (1) Nabilone [±]; and

15 (2) Dronabinol (-)-delta-9-trans tetrahydrocannabinol in  
16 an oral solution in a drug product approved for  
17 marketing by the United States Food and Drug  
18 Administration."

19 SECTION 3. Statutory material to be repealed is bracketed  
20 and stricken. New statutory material is underscored.

21 SECTION 4. This Act shall take effect upon its approval.



H.B. NO. 2385  
H.D. 2  
S.D. 1

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APPROVED this 10 day of JUL , 2018



GOVERNOR OF THE STATE OF HAWAII

HB No. 2385, HD 2, SD 1

THE HOUSE OF REPRESENTATIVES OF THE STATE OF HAWAII

Date: April 27, 2018  
Honolulu, Hawaii

We hereby certify that the above-referenced Bill on this day passed Final Reading in the House of Representatives of the Twenty-Ninth Legislature of the State of Hawaii, Regular Session of 2018.



Scott K. Saiki  
Speaker  
House of Representatives




Brian L. Takeshita  
Chief Clerk  
House of Representatives


H.B. No. 2385, H.D. 2, S.D. 1

**THE SENATE OF THE STATE OF HAWAI'I**

Date: April 10, 2018  
Honolulu, Hawai'i 96813

We hereby certify that the foregoing Bill this day passed Third Reading in the  
Senate of the Twenty-ninth Legislature of the State of Hawai'i, Regular Session of 2018.

  
President of the Senate

  
Clerk of the Senate