
A BILL FOR AN ACT

RELATING TO CONTROLLED SUBSTANCES.

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 subsection (g) to read as follows:
3 "(g) Any of the following cannabinoids, their salts,
4 isomers, and salts of isomers, unless specifically excepted,
5 whenever the existence of these salts, isomers, and salts of
6 isomers is possible within the specific chemical designation:
7 (1) Tetrahydrocannabinols; meaning tetrahydrocannabinols
8 naturally contained in a plant of the genus Cannabis
9 (cannabis plant), as well as synthetic equivalents of
10 the substances contained in the plant, or in the
11 resinous extractives of Cannabis, sp. or synthetic
12 substances, derivatives, and their isomers with
13 similar chemical structure and pharmacological
14 activity to those substances contained in the plant,
15 such as the following: Delta 1 cis or trans
16 tetrahydrocannabinol, and their optical isomers; Delta
17 6 cis or trans tetrahydrocannabinol [7] (Other name:



- 1 Delta 8 tetrahydrocannabinol), and their optical
2 isomers; and Delta 3,4 cis or trans-
3 tetrahydrocannabinol, and its optical isomers (since
4 nomenclature of these substances is not
5 internationally standardized, compounds of these
6 structures, regardless of numerical designation of
7 atomic positions, are covered);
- 8 (2) Naphthoylindoles; meaning any compound containing a
9 3-(1-naphthoyl)indole structure with substitution at
10 the nitrogen atom of the indole ring by a alkyl,
11 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2-piperidinyl)methyl or
13 2-(4-morpholinyl)ethyl group, whether or not further
14 substituted in the indole ring to any extent and
15 whether or not substituted in the naphthyl ring to any
16 extent;
- 17 (3) Naphthylmethylinindoles; meaning any compound containing
18 a 1H-indol-3-yl-(1-naphthyl) methane structure with
19 substitution at the nitrogen atom of the indole ring
20 by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
21 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or



- 1 2-(4-morpholinyl) ethyl group whether or not further
2 substituted in the indole ring to any extent and
3 whether or not substituted in the naphthyl ring to any
4 extent;
- 5 (4) Naphthoylpyrroles; meaning any compound containing a
6 3-(1-naphthoyl)pyrrole structure with substitution at
7 the nitrogen atom of the pyrrole ring by a alkyl,
8 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
9 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)
10 ethyl group whether or not further substituted in the
11 pyrrole ring to any extent, whether or not substituted
12 in the naphthyl ring to any extent;
- 13 (5) Naphthylmethylindenes; meaning any compound containing
14 a naphthylideneindene structure with substitution at
15 the 3-position of the indene ring by a alkyl,
16 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
17 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)
18 ethyl group whether or not further substituted in the
19 indene ring to any extent, whether or not substituted
20 in the naphthyl ring to any extent;



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



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



- 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)-1 H-indazole-3-carbonyl)-L-
12 valinate, and geometric isomers, salts, and salts of
13 isomers (Other names: FUB-AMB);
- 14 (21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
15 carboxamido)-3-methylbutanoate, and geometric isomers,
16 salts, and salts of isomers (Other names: 5-fluoro-
17 AMB, 5-fluoro-AMP);
- 18 (22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-
19 indazole-3-carboxamide, and geometric isomers, salts,
20 and salts of isomers (Other names: AKB48 N-(5-



- 1 fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
2 analog, 5F-APINACA);
- 3 (23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
4 geometric isomers, salts, and salts of isomers (Other
5 names: STS-135, 5F-APICA; 5-fluoro-APICA);
- 6 (24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
7 carboxylate, and geometric isomers, salts, and salts
8 of isomers (Other names: NM2201);
- 9 (25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
10 (cyclohexylmethyl)-1H-indazole-3-carboxamide, and
11 geometric isomers, salts, and salts of isomers (Other
12 names: MAB-CHMINACA and ADB-CHMINACA);
- 13 (26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
14 carboxamido]-3,3-dimethylbutanoate (Other names: 5F-
15 ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical,
16 positional, and geometric isomers, salts, and salts of
17 isomers; and
- 18 (27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)indazole-3-
19 carboxamide (CUMYL-4CN-BINACA), its optical,
20 positional, and geometric isomers, salts, and salts of
21 isomers; also known as SGT-78, 4-CN-CUMYL-BINACA;



H.B. NO. 422

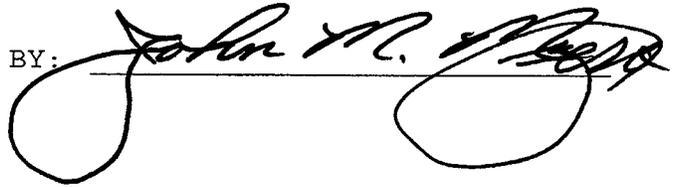
1 CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano CUMYL-
2 BUTINACA."

3 SECTION 2. Statutory material to be repealed is bracketed
4 and stricken. New statutory material is underscored.

5 SECTION 3. This Act shall take effect upon its approval.

6

INTRODUCED BY:



JAN 22 2021



H.B. NO. 422

Report Title:

Uniform Controlled Substances Act; Schedule I; Cannabinoids;
Tetrahydrocannabinols

Description:

Adds Delta 8 tetrahydrocannabinol as another name for Delta 6
cis or trans tetrahydrocannabinol, on schedule I of the list of
controlled substances.

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

