
A BILL FOR AN ACT

RELATING TO CONTROLLED SUBSTANCES.

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

1 SECTION 1. Section 328G-3, Hawaii Revised Statutes, is
2 amended by amending subsection (f) to read as follows:

3 "(f) No person shall sell, hold, offer, or distribute for
4 sale any hemp product into which a synthetic cannabinoid has
5 been added[-] or that contains cannabinoids created through
6 isomerization, including but not limited to Delta 8
7 tetrahydrocannabinol."

8 SECTION 2. Section 329-14, Hawaii Revised Statutes, is
9 amended by amending subsection (g) to read as follows:

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

14 (1) Tetrahydrocannabinols; meaning tetrahydrocannabinols
15 naturally contained in a plant of the genus Cannabis
16 (cannabis plant), as well as synthetic equivalents of
17 the substances contained in the plant, or in the



1 resinous extractives of Cannabis, sp. or synthetic
2 substances, derivatives, and their isomers with
3 similar chemical structure and pharmacological
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; Delta 8 cis or trans tetrahydrocannabinol,
9 and their optical isomers; and Delta 3,4 cis or trans-
10 tetrahydrocannabinol, and its optical isomers (since
11 nomenclature of these substances is not
12 internationally standardized, compounds of these
13 structures, regardless of numerical designation of
14 atomic positions, are covered);

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



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

3 (3) Naphthylmethyloindoles; meaning any compound containing
4 a 1H-indol-3-yl-(1-naphthyl) methane 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 and
10 whether or not substituted in the naphthyl ring to any
11 extent;

12 (4) Naphthoylpyrroles; meaning any compound containing a
13 3-(1-naphthoyl)pyrrole structure with substitution at
14 the nitrogen atom of the pyrrole 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 pyrrole ring to any extent, whether or not substituted
19 in the naphthyl ring to any extent;

20 (5) Naphthylmethyloindenes; meaning any compound containing
21 a naphthylideneindene structure with substitution at



1 the 3-position of the indene ring by a alkyl,
2 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
3 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)
4 ethyl group whether or not further substituted in the
5 indene ring to any extent, whether or not substituted
6 in the naphthyl ring to any extent;

7 (6) Phenylacetylindoles; meaning any compound containing a
8 3-phenylacetylindole structure with substitution at
9 the nitrogen atom of the indole ring by a alkyl,
10 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
11 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)
12 ethyl group whether or not further substituted in the
13 indole ring to any extent, whether or not substituted
14 in the phenyl ring to any extent;

15 (7) Cyclohexylphenols; meaning any compound containing a
16 2-(3-hydroxycyclohexyl) phenol structure with
17 substitution at the 5-position of the phenolic ring by
18 a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
19 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
20 2-(4-morpholinyl) ethyl group whether or not
21 substituted in the cyclohexyl ring to any extent;



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



1 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
2 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,
3 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
4 morpholinyl)methyl, or tetrahydropyranylmethyl group,
5 whether or not further substituted in the indole ring
6 to any extent and whether or not substituted in the
7 tetramethylcyclopropyl ring to any extent;

8 (12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
9 its optical, positional, and geometric isomers, salts,
10 and salts of isomers (Other names: APINACA, AKB48);

11 (13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
12 optical, positional, and geometric isomers, salts, and
13 salts of isomers (Other names: PB-22; QUPIC);

14 (14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-
15 carboxylate, its optical, positional, and geometric
16 isomers, salts, and salts of isomers (Other names: 5-
17 fluoro-PB-22; 5F-PB-22);

18 (15) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-
19 fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
20 positional, and geometric isomers, salts, and salts of
21 isomers (Other names: AB-FUBINACA);



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



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



1 positional, and geometric isomers, salts, and salts of
2 isomers;

3 (27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-
4 3-carboxamide, its optical, positional, and geometric
5 isomers, salts, and salts of isomers (Other names:
6 SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-
7 CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA;
8 CUMYL-4CN-BINACA);

9 (28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
10 fluoropentyl)-1H-indazole-3-carboxamide (Other name:
11 5F-AB-PINACA);

12 (29) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
13 carboxamido)-3-methylbutanoate (Other names: MMB-
14 CHMICA; AMB-CHMICA);

15 (30) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
16 pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-
17 CUMYL-P7AICA); and

18 (31) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
19 3-carboxamido)butanoate (MDMB-4en-PINACA)."

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

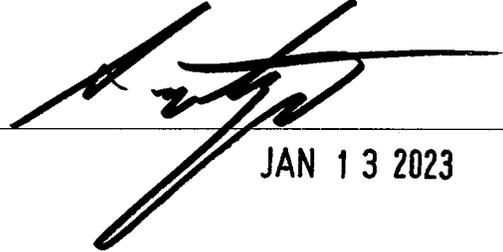


H.B. NO. 10

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

2

INTRODUCED BY:



JAN 13 2023



H.B. NO. 70

Report Title:

Hemp Products; Uniform Controlled Substances Act; Schedule I;
Cannabinoids; Delta 8 Tetrahydrocannabinol

Description:

Bans hemp products that contain Delta 8 tetrahydrocannabinol.
Adds Delta 8 tetrahydrocannabinol to the list of schedule I
controlled substances.

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