
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 (g) to read as follows:

3 "(g) No person shall sell~~[-]~~ or hold, offer, or distribute
4 for sale any crude extract or manufactured hemp product [~~into~~]:

5 (1) Into which an artificially derived cannabinoid or a
6 synthetic cannabinoid has been added~~[-]~~; or

7 (2) That contains cannabinoids created through
8 isomerization, including Delta 6 cis or trans
9 tetrahydrocannabinol, and their optical isomers (other
10 names: Delta 8 cis or trans tetrahydrocannabinol, and
11 their optical isomers)."

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

14 "(g) Cannabinoids. Unless specifically excepted or unless
15 listed in another schedule, any of the following cannabinoids,
16 including their salts, isomers, and salts of isomers, whenever



1 the existence of these salts, isomers, and salts of isomers is
2 possible within the specific chemical designation:

3 (1) Tetrahydrocannabinols; meaning tetrahydrocannabinols
4 naturally contained in a plant of the genus Cannabis
5 (cannabis plant), as well as synthetic equivalents of
6 the substances contained in the plant, or in the
7 resinous extractives of Cannabis, sp. or synthetic
8 substances, derivatives, and their isomers with
9 similar chemical structure and pharmacological
10 activity to those substances contained in the plant,
11 such as the following: Delta 1 cis or trans
12 tetrahydrocannabinol, and their optical isomers; Delta
13 6 cis or trans tetrahydrocannabinol, and their optical
14 isomers[+] (other names: Delta 8 cis or trans
15 tetrahydrocannabinol, and their optical isomers); and
16 Delta 3,4 cis or trans-tetrahydrocannabinol, and its
17 optical isomers (since nomenclature of these
18 substances is not internationally standardized,
19 compounds of these structures, regardless of numerical
20 designation of atomic positions, are covered);



- 1 (2) Naphthoylindoles; meaning any compound containing a 3-
2 (1-naphthoyl)indole structure with substitution at the
3 nitrogen atom of the indole ring by [a] an 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 naphthyl ring to any
9 extent;
- 10 (3) Naphthylmethylinindoles; meaning any compound containing
11 a 1H-indol-3-yl-(1-naphthyl) methane structure with
12 substitution at the nitrogen atom of the indole ring
13 by [a] an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
14 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
15 2-(4-morpholinyl) ethyl group whether or not further
16 substituted in the indole ring to any extent and
17 whether or not substituted in the naphthyl ring to any
18 extent;
- 19 (4) Naphthoylpyrroles; meaning any compound containing a
20 3-(1-naphthoyl)pyrrole structure with substitution at
21 the nitrogen atom of the pyrrole ring by [a] an alkyl,



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



- 1 (7) Cyclohexylphenols; meaning any compound containing a
2 2-(3-hydroxycyclohexyl) phenol structure with
3 substitution at the 5-position of the phenolic ring by
4 [a] an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
5 cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
6 2-(4-morpholinyl) ethyl group whether or not
7 substituted in the cyclohexyl ring to any extent;
- 8 (8) Benzoylindoles; meaning any compound containing a 3-
9 (benzoyl) indole structure with substitution at the
10 nitrogen atom of the indole ring by [a] an alkyl,
11 haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2-piperidinyl) methyl, or 2-(4-
13 morpholinyl) ethyl group whether or not further
14 substituted in the indole ring to any extent and
15 whether or not substituted in the phenyl ring to any
16 extent;
- 17 (9) [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
18 pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-
19 naphthalenylmethanone (another trade name is WIN
20 55,212-2);



- 1 (10) (6a,10a)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-
2 methyloctan-2-yl)-6a,7,10,10a-
3 tetrahydrobenzo[c]chromen-1-ol (Other trade names are:
4 HU-210/HU-211);
- 5 (11) Tetramethylcyclopropanoylindoles; meaning any compound
6 containing a 3-tetramethylcyclopropanoylindole
7 structure with substitution at the nitrogen atom of
8 the indole ring by an alkyl, haloalkyl, cyanoalkyl,
9 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
10 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,
11 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
12 morpholinyl)methyl, or tetrahydropyranylmethyl group,
13 whether or not further substituted in the indole ring
14 to any extent and whether or not substituted in the
15 tetramethylcyclopropyl ring to any extent;
- 16 (12) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
17 its optical, positional, and geometric isomers, salts,
18 and salts of isomers (Other names: APINACA, AKB48);
- 19 (13) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
20 optical, positional, and geometric isomers, salts, and
21 salts of isomers (Other names: PB-22; QUPIC);



- 1 (14) Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-
2 carboxylate, its optical, positional, and geometric
3 isomers, salts, and salts of isomers (Other names: 5-
4 fluoro-PB-22; 5F-PB-22);
- 5 (15) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-
6 fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
7 positional, and geometric isomers, salts, and salts of
8 isomers (Other names: AB-FUBINACA);
- 9 (16) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
10 indazole-3-carboxamide, its optical, positional, and
11 geometric isomers, salts, and salts of isomers (Other
12 names: ADB-PINACA);
- 13 (17) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
14 (cyclohexylmethyl)-1H-indazole-3-carboxamide, its
15 optical, positional, and geometric isomers, salts, and
16 salts of isomers (Other names: AB-CHMINACA);
- 17 (18) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
18 indazole-3-carboxamide, and geometric isomers, salts,
19 and salts of isomers (Other names: AB-PINACA);



- 1 (19) [1-(5-fluoropentyl)-1H-indazol-3-yl] (naphthalen-1-
2 yl)methanone, and geometric isomers, salts, and salts
3 of isomers (Other names: THJ-2201);
- 4 (20) Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-
5 valinate, and geometric isomers, salts, and salts of
6 isomers (Other names: FUB-AMB, Methyl 2-(1-(4-
7 fluorobenzyl)-1H-indazole-3-carboxamido)-3-
8 methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);
- 9 (21) (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
10 carboxamido)-3-methylbutanoate, and geometric isomers,
11 salts, and salts of isomers (Other names: 5-fluoro-
12 AMB, 5-fluoro-AMP);
- 13 (22) N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-
14 indazole-3-carboxamide, and geometric isomers, salts,
15 and salts of isomers (Other names: AKB48 N-(5-
16 fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
17 analog, 5F-APINACA);
- 18 (23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
19 geometric isomers, salts, and salts of isomers (Other
20 names: STS-135, 5F-APICA; 5-fluoro-APICA);



- 1 (24) Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
2 carboxylate, and geometric isomers, salts, and salts
3 of isomers (Other names: NM2201; CBL2201);
- 4 (25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
5 (cyclohexylmethyl)-1H-indazole-3-carboxamide, and
6 geometric isomers, salts, and salts of isomers (Other
7 names: MAB-CHMINACA and ADB-CHMINACA);
- 8 (26) Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
9 carboxamido]-3,3-dimethylbutanoate (Other names: 5F-
10 ADB, [~~5-fluoro-ADB~~,] 5-fluoro-ADB, and 5F-MDMB-
11 PINACA), its optical, positional, and geometric
12 isomers, salts, and salts of isomers;
- 13 (27) 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-
14 3-carboxamide, its optical, positional, and geometric
15 isomers, salts, and salts of isomers (Other names:
16 SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA; CUMYL-
17 CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-BUTINACA;
18 CUMYL-4CN-BINACA);
- 19 (28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
20 fluoropentyl)-1H-indazole-3-carboxamide (Other name:
21 5F-AB-PINACA);



- 1 (29) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
2 carboxamido)-3-methylbutanoate (Other names: MMB-
3 CHMICA; AMB-CHMICA);
- 4 (30) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
5 pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-
6 CUMYL-P7AICA);
- 7 (31) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
8 3-carboxamido)butanoate (MDMB-4en-PINACA);
- 9 (32) Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
10 carboxamido)-3,3-dimethylbutanoate (Other name: 5F-
11 EDMB-PINACA);
- 12 (33) Methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-
13 3,3-dimethylbutanoate (Other names: 5F-MDMB-PICA; 5F-
14 MDMB-2201);
- 15 (34) N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-
16 carboxamide (Other names: FUB-AKB48; FUB-APINACA;
17 AKB48 N-(4-FLUOROBENZYL));
- 18 (35) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
19 indazole-3-carboxamide (Other names: 5F-CUMYL-PINACA;
20 SGT-25); and



1 (36) (1-(4-fluorobenzyl)-1H-indol-3-yl) (2,2,3,3-
2 tetramethylcyclopropyl)methanone (Other name: FUB-
3 144) "

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

6 SECTION 4. This Act shall take effect on December 31,
7 2050.



Report Title:

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

Description:

Prohibits the sale or holding, offering, and distribution of hemp products that contain cannabinoids created through isomerization. Inserts the alternative name for Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers into the list of schedule I controlled substances. Takes effect 12/31/2050. (SD1)

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

