A BILL FOR AN ACT

RELATING TO CONTROLLED SUBSTANCES.

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

1	SECT	ION 1. Section 328G-3, Hawaii Revised Statutes, is
2	amended b	y amending subsection (f) to read as follows:
3	"(f)	No person shall sell[τ] or hold, offer, or distribute
4	for sale	any hemp product [into]:
5 .	(1)	Into which a synthetic cannabinoid has been added[-];
6		<u>or</u>
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); provided that this paragraph
12		shall not be construed to prohibit a medical cannabis
13		dispensary licensed pursuant to chapter 329D from
14		selling or holding, offering, or distributing for sale
15		cannabis or manufactured cannabis products that
16		contain naturally-occurring Delta 8
17		tetrahydrocannabinol."

1	SECTION 2. 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, and their optical
18	isomers[+] (other names: Delta 8 cis or trans
19	tetrahydrocannabinol, and their optical isomers); and
20	Delta 3,4 cis or trans-tetrahydrocannabinol, and its
21	optical isomers (since nomenclature of these

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

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

1		ethyl group whether or not further substituted in the
2		indole ring to any extent, whether or not substituted
3		in the phenyl ring to any extent;
4	(7)	Cyclohexylphenols; meaning any compound containing a
5		2-(3-hydroxycyclohexyl) phenol structure with
6		substitution at the 5-position of the phenolic ring by
7		a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
8		cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
9		2-(4-morpholinyl) ethyl group whether or not
10		substituted in the cyclohexyl ring to any extent;
11	(8)	Benzoylindoles; meaning any compound containing a
12		3-(benzoyl) indole structure with substitution at the
13		nitrogen atom of the indole ring by a alkyl,
14		haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
15		1-(N-methyl-2-piperidinyl) methyl, or 2-(4-
16		morpholinyl) ethyl group whether or not further
17		substituted in the indole ring to any extent and
18		whether or not substituted in the phenyl ring to any
19		extent;
20	(9)	[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
21		pvrrolo[1.2.3-de]-1. 4-benzoxazin-6-v]]-1-

H.B. NO. H.D. 1

```
1
              naphthalenylmethanone (another trade name is WIN
2
              55,212-2);
3
        (10)
              (6a, 10a) - 9 - (hydroxymethy1) - 6, 6 - dimethy1 - 3 - (2 - 6)
4
              methyloctan-2-yl)-6a,7,10,10a-
5
              tetrahydrobenzo[c]chromen-1-ol (Other trade names are:
6
              HU-210/HU-211);
7
        (11)
              Tetramethylcyclopropanoylindoles; meaning any compound
8
              containing a 3-tetramethylcyclopropanoylindole
9
              structure with substitution at the nitrogen atom of
10
              the indole ring by an alkyl, haloalkyl, cyanoalkyl,
11
              alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
12
              methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl,
13
              1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
14
              morpholinyl) methyl, or tetrahydropyranyl methyl group,
15
              whether or not further substituted in the indole ring
16
              to any extent and whether or not substituted in the
17
              tetramethylcyclopropyl ring to any extent;
18
        (12)
              N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
19
              its optical, positional, and geometric isomers, salts,
20
              and salts of isomers (Other names: APINACA, AKB48);
```

H.B. NO. H.D. 1

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

```
1
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
        (18)
2
              indazole-3-carboxamide, and geometric isomers, salts,
3
              and salts of isomers (Other names: AB-PINACA);
4
              [1-(5-fluoropentyl)-1H-indazol-3-vl](naphthalen-1-
        (19)
5
              yl) methanone, and geometric isomers, salts, and salts
6
              of isomers (Other names: THJ-2201);
7
        (20)
              Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-
8
              valinate, and geometric isomers, salts, and salts of
9
              isomers (Other names: FUB-AMB, Methyl 2-(1-(4-
10
              fluorobenzyl)-1H-indazole-3-carboxamido)-3-
11
              methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);
12
        (21)
              (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
13
              carboxamido)-3-methylbutanoate, and geometric isomers,
14
              salts, and salts of isomers (Other names: 5-fluoro-
15
              AMB, 5-fluoro-AMP);
16
        (22)
              N-((3s,5s,7s)-adamantan-1-y1)-1-(5-fluoropenty1)-1H-
17
              indazole-3-carboxamide, and geometric isomers, salts,
18
              and salts of isomers (Other names: AKB48 N-
19
              (5-fluoropentyl) analog, 5F-AKB48, APINACA 5-
20
              fluoropentyl analog, 5F-APINACA);
```

```
1
        (23) N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
2
              geometric isomers, salts, and salts of isomers (Other
3
              names: STS-135, 5F-APICA; 5-fluoro-APICA);
4
              Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
        (24)
5
              carboxylate, and geometric isomers, salts, and salts
6
              of isomers (Other names: NM2201; CBL2201);
7
        (25)
              N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
8
              (cyclohexylmethyl)-1H-indazole-3-carboxamide, and
9
              geometric isomers, salts, and salts of isomers (Other
10
              names: MAB-CHMINACA and ADB-CHMINACA);
11
        (26)
              Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
12
              carboxamido]-3,3-dimethylbutanoate (Other names:
13
              5F-ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its
14
              optical, positional, and geometric isomers, salts, and
              salts of isomers;
15
16
        (27)
              1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-
17
              3-carboxamide, its optical, positional, and geometric
18
              isomers, salts, and salts of isomers (Other names:
19
              SGT-78; 4-CN-CUMYL BINACA; 4-CN-CUMYL-BUTINACA;
20
              CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano-CUMYL-
21
              BUTINACA; CUMYL-4CN-BINACA);
```

```
1
        (28) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
2
              fluoropentyl)-1H-indazole-3-carboxamide (Other name:
3
              5F-AB-PINACA);
4
        (29)
              Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
5
              carboxamido)-3-methylbutanoate (Other names:
6
              MMB-CHMICA; AMB-CHMICA);
7
        (30)
              1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
8
              pyrrolo[2,3-b]pyridine-3-carboxamide (Other names:
9
              5F-CUMYL-P7AICA); and
10
        (31)
              Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
11
              3-carboxamido)butanoate (MDMB-4en-PINACA)."
12
         SECTION 3. Statutory material to be repealed is bracketed
13
    and stricken. New statutory material is underscored.
14
         SECTION 4. This Act shall take effect on June 30, 3000.
```

Report Title:

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

Description:

Prohibits the sale and distribution of hemp products that contain cannabinoids created through isomerization, including Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers (other names: Delta 8 cis or trans tetrahydrocannabinol, and their optical isomers). Exempts certain cannabis and manufactured cannabis product sales from the prohibition. Inserts alternative names for Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers into the list of schedule I controlled substances. Effective 6/30/3000. (HB70 HD1)

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