## 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 as follows: 3 1. By amending subsection (b) to read: 4 "(b) Any of the following opiates, including their 5 isomers, esters, ethers, salts, and salts of isomers, esters, 6 and ethers, unless specifically excepted, whenever the existence 7 of these isomers, esters, ethers, and salts is possible within 8 the specific chemical designation: 9 (1)Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-10 phenethyl) -4-piperidinyl] -N-phenylacetamide); 11 (2) Acetylmethadol; 12 (3) Allylprodine; 13 (4)Alphacetylmethadol (except levo-alphacetylmethadol, 14 levomethadyl acetate, or LAAM); **15** (5) Alphameprodine; 16 (6) Alphamethadol;

```
1
          (7)
               Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-
 2
               phenyl)ethyl-4-piperidyl] propionanilide; 1-(1-methyl-
 3
               2-phenylethyl)-4-(N-propanilido) piperidine);
 4
          (8)
               Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
 5
               thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
 6
          (9)
               Benzethidine;
7
         (10)
               Betacetylmethadol;
8
         (11)
               Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-
9
               piperidinyl] -N-phenylpropanamide);
10
         (12)
               Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-
11
               phenethyl)-3-methyl-4-piperidinyl]-N-
12
               phenylpropanamide);
13
         (13)
               Betameprodine;
14
         (14)
               Betamethadol;
15
        (15)
               Betaprodine;
16
               Clonitazene;
        (16)
17
        (17)
               Dextromoramide;
18
        (18)
               Diampromide;
19
        (19)
               Diethylthiambutene;
20
        (20)
               Difenoxin;
21
               Dimenoxadol;
        (21)
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1
         (22)
               Dimepheptanol;
 2
         (23)
               Dimethylthiambutene;
 3
         (24)
               Dioxaphetyl butyrate;
 4
         (25)
               Dipipanone;
 5
         (26)
               Ethylmethylthiambutene;
 6
         (27)
               Etonitazene;
 7
         (28)
               Etoxeridine;
 8
         (29)
               Furethidine;
 9
         (30)
               Hydroxypethidine;
10
         (31)
               Ketobemidone;
11
         (32)
               Levomoramide;
12
         (33)
               Levophenacylmorphan;
13
               3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-
         (34)
14
               piperidyl]-N-phenylpropanamide);
15
               3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-
         (35)
16
               4-piperidinyl]-N-phenylpropanamide);
17
         (36)
               Morpheridine;
18
         (37)
               MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
19
         (38)
               Noracymethadol;
20
         (39)
               Norlevorphanol;
21
         (40)
               Normethadone;
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1
        (41)
               Norpipanone;
2
        (42)
               Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-
3
               phenethyl)-4-piperidinyl] propanamide;
4
        (43)
               PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine;
5
        (44)
               Phenadoxone;
        (45)
6
               Phenampromide;
7
        (46)
               Phenomorphan;
8
        (47)
               Phenoperidine;
9
        (48)
               Piritramide;
10
        (49) Proheptazine;
11
        (50)
               Properidine;
12
        (51)
               Propiram;
13
        (52)
               Racemoramide;
14
               Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
        (53)
15
               piperidinyl]-propanamide);
16
        (54)
               Tilidine;
17
               Trimeperidine;
        (55)
18
         (56)
               N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
19
               (benzylfentanyl), its optical isomers, salts, and
20
               salts of isomers;
```

1	(57)	N-[1-(2-thienyl)methyl-4-piperidyl]-N-
2		phenylpropanamide (thenylfentanyl), its optical
3		isomers, salts, and salts of isomers; [and]
4	(58)	N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide,
5		(acetyl fentanyl), its optical, positional, and
6		geometric isomers, salts, and salts of isomers $[-]$ :
7	(59)	AH-7921 (3,4-dichloro-N-[(1- dimethylamino)
8		cyclohexylmethyl]benzamide), its isomers, esters,
9		ethers, salts, and salts of isomers, esters, and
10		ethers;
11	(60)	N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, its
12		isomers, esters, ethers, salts, and salts of isomers,
13		esters, and ethers (Other names: Butyryl fentanyl);
14	(61)	N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-
15		yl]-N-phenylpropionamide, its isomers, esters, ethers,
16		salts, and salts of isomers, esters, and ethers (Other
17	e e	<pre>names: beta-hydroxythiofentanyl);</pre>
18	(62)	N-(1-phenthylpiperidin-4-y1)-N-phenylfuran-2-
19		carboxamide, its isomers, esters, ethers, salts, and
20		salts of isomers, esters, and ethers (other names:
21		Furanyl fentanyl); and

1	(63)	3,4-dicholoro- $N$ -[2-(dimethylamino)cyclohexyl]- $N$ -
2		methylbenzamide, its isomers, esters, ethers, salts,
3		and salts of isomers, esters, and ethers (Other names:
4		<u>U-47700).</u> "
5	2.	By amending subsection (g) to read:
6	<b>"</b> (g)	Any of the following cannabinoids, their salts,
7	isomers,	and salts of isomers, unless specifically excepted,
8	whenever	the existence of these salts, isomers, and salts of
9	isomers i	s possible within the specific chemical designation:
10	(1)	Tetrahydrocannabinols; meaning tetrahydrocannabinols
11		naturally contained in a plant of the genus Cannabis
12		(cannabis plant), as well as synthetic equivalents of
13		the substances contained in the plant, or in the
14		resinous extractives of Cannabis, sp. or synthetic
15		substances, derivatives, and their isomers with
16		similar chemical structure and pharmacological
17		activity to those substances contained in the plant,
18		such as the following: Delta 1 cis or trans
19		tetrahydrocannabinol, and their optical isomers; Delta
20		6 cis or trans tetrahydrocannabinol, and their optical
21		isomers; and Delta 3,4 cis or trans-

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

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

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1
          (9)
               2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
2
              pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-
3
              napthalenylmethanone (another trade name is WIN
4
               55,212-2);
5
        (10)
              (6a, 10a) - 9 - (hydroxymethyl) - 6, 6 - dimethyl - 3 - (2 - a)
6
              methyloctan-2-yl)-6a,7,10,10a-
7
              tetrahydrobenzo[c]chromen-1-ol (other trade names are:
8
              HU-210/HU-211);
9
        (11)
              Tetramethylcyclopropanoylindoles; meaning any compound
10
              containing a 3-tetramethylcyclopropanoylindole
11
              structure with substitution at the nitrogen atom of
12
              the indole ring by an alkyl, haloalkyl, cyanoalkyl,
              alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
13
14
              methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl,
15
              1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
              morpholinyl) methyl, or tetrahydropyranylmethyl group,
16
17
              whether or not further substituted in the indole ring
18
              to any extent and whether or not substituted in the
19
              tetramethylcyclopropyl ring to any extent[-];
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# S.B. NO. 997 S.D. 1

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1
         (12)
               N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
2
               its optical, positional, and geometric isomers, salts,
               and salts of isomers. (Other names: APINACA, AKB48);
3
4
               Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
         (13)
5
               optical, positional, and geometric isomers, salts, and
6
             salts of isomers (Other names: PB-22; OUPIC);
7
               Quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-
         (14)
8
               carboxylate, its optical, positional, and geometric
9
               isomers, salts, and salts of isomers (Other names: 5-
10
               fluoro-PB-22; 5F-PB-22);
11
               N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-amino-3-methyl-1-oxobutan-2-yl)
         (15)
12
               fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
13
               positional, and geometric isomers, salts, and salts of
14
               isomers (Other names: AB-FUBINACA);
15
        (16)
               N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
16
               indazole-3-carboxamide, its optical, positional, and
17
               geometric isomers, salts, and salts of isomers (Other
18
               names: ADB-PINACA);
19
        (17)
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
20
               (cyclohexylmethyl) -1H-indazole-3-carboxamide, its
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1
              optical, positional, and geometric isomers, salts, and
2
              salts of isomers (Other names: AB-CHMINACA);
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
3
        (18)
              indazole-3-carboxamide, and geometric isomers, salts,
4
5
              and salts of isomers (Other names: AB-PINACA);
              [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-
6
        (19)
7
              yl) methanone, and geometric isomers, salts, and salts
8
              of isomers (Other names: THJ-2201);
9
              Methyl (1-(4-fluorobenzyl)-1 H-indazole-3-carbonyl)-L-
        (20)
10
              valinate, and geometric isomers, salts, and salts of
11
              isomers (Other names: FUB-AMB);
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
              N-(3s, 5s, 7s) -adamantan-1-yl) -1-(5-fluoropentyl) -1H-
        (22)
17
              indazole-3-carboxamide, and geometric isomers, salts,
18
              and salts of isomers (Other names: AKB48 N-(5-
              fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
19
              analog, 5F-APINACA);
20
```

```
1
              N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
        (23)
 2
              geometric isomers, salts, and salts of isomers (Other
              names: STS-135, 5F-APICA; 5-fluoro-APICA);
 3
 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); [and]
7
              N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
        (25)
 8
               (cyclohexylmethyl) -1H-indazole-3-carboxamide, and
 9
              geometric isomers, salts, and salts of isomers (Other
10
              names: MAB-CHMINACA and ADB-CHMINACA) [-]; and
11
              Methyl -2-[1-(5-fluoropentyl)-1H-indazole-3-
        (26)
12
              carboxamido]-3,3-dimethylbutanoate (other names: 5F-
              ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical,
13
14
              positional, and geometric isomers, salts, and salts of
15
              isomers."
16
         SECTION 2. Section 329-16, Hawaii Revised Statutes, is
17
    amended by amending subsection (c) to read as follows:
18
         "(c) Any of the following opiates, including their
19
    isomers, esters, ethers, salts, and salts of isomers, whenever
20
    the existence of these isomers, esters, ethers, and salts is
21
    possible within the specific chemical designation:
```



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1
         (1) Alfentanil;
2
         (2) Alphaprodine;
3
         (3) Anileridine;
4
         (4)
             Bezitramide;
5
         (5)
              Bulk Dextropropoxyphene (nondosage form);
6
         (6)
              Carfentanil:
7
         (7)
             Dihydrocodeine;
8
         (8)
              Diphenoxylate;
9
         (9)
              Fentanyl;
10
              Isomethadone;
        (10)
11
       (11)
              Levo-alphacetylmethadol (LAAM);
12
        (12)
              Levomethorphan;
13
              Levorphanol;
        (13)
14
        (14)
              Metazocine;
15
        (15)
             Methadone;
16
              Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-
        (16)
17
              diphenyl butane;
18
        (17)
              Moramide-Intermediate, 2-methyl-3-morpholino-1, 1
19
              diphenyl-propane-carboxylic acid;
20
              Pethidine (Meperidine);
        (18)
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1
        (19)
              Pethidine-Intermediate-A, 4-cyano-1-methyl-4-
2
              phenylpiperidine;
3
        (20)
              Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-
4
              carboxylate;
5
        (21)
              Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-
6
              4-carboxylic acid;
7
        (22)
              Phenazocine;
8
        (23)
              Piminodine;
9
        (24)
              Racemethorphan;
10
        (25)
              Racemorphan;
11
              Remifentanil;
        (26)
12
        (27)
              Sufentanil; [and]
13
        (28)
              Tapentadol [-]; and
14
        (29)
              Thiafentanil."
         SECTION 3. Section 329-22, Hawaii Revised Statutes, is
15
16
    amended by amending subsection (d) to read as follows:
17
               Depressants. Unless specifically exempted or
18
    excluded or unless listed in another schedule, any material,
19
    compound, mixture, or preparation that contains any quantity of
20
    the following substances having a depressant effect on the
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1
    central nervous system, including its salts, isomers, and salts
 2
    of isomers:
 3
         (1) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-
 4
              propionamide], (Vimpat); [and]
 5
         (2) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic
 6
              acid][-]; and
 7
         (3) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-
 8
              yl]butanamide) (Other names: BRV; UCB-34714; Briviact)
9
              and its salts."
10
         SECTION 4. Statutory material to be repealed is bracketed
11
    and stricken. New statutory material is underscored.
12
         SECTION 5. This Act shall take effect on January 7, 2059.
13
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### Report Title:

Uniform Controlled Substances Act

#### Description:

Updates the Uniform Controlled Substances Act to make it consistent with amendments in federal controlled substances law as required under the authority to schedule controlled substances. Takes effect on 1/7/2059. (SD1)

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