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: "(b) Any of the following opiates, including their 4 isomers, esters, ethers, salts, and salts of isomers, esters, 5 6 and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers, and salts is possible within 7 8 the specific chemical designation: 9 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-(1)10 phenethyl) -4-piperidinyl]-N-phenylacetamide); 11 Acetylmethadol; (2) 12 (3) Acetyl fentanyl (N-(1-phenethylpiperidin-4-yl)-Nphenylacetamide), its optical, positional, and 13 14 geometric isomers, salts, and salts of isomers; 15 (4) Acryl fentanyl [N-(1-phenethylpiperidin-4-yl)-N-16 phenylacrylamide] (Other name: acryloylfentanyl);

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1
          (5) AH-7921 (3,4-dichloro-N-[(1-
2
               dimethylamino) cyclohexylmethyl]benzamide);
3
         [\frac{(3)}{(3)}] (6) Allylprodine;
4
         [<del>(4)</del>] (7) Alphacetylmethadol (except levo-
5
               alphacetylmethadol, levomethadyl acetate, or LAAM);
6
         [\frac{(5)}{(8)}] (8) Alphameprodine;
7
         [(6)] (9) Alphamethadol;
8
         [\frac{(7)}{1}] (10) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-
9
               phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-
10
               2-phenylethyl)-4-(N-propanilido)piperidine);
11
         [+(8)] (11) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
12
               thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
13
         [+9+] (12) Benzethidine;
14
         (13) Benzylfentanyl (N-[1-benzyl-4-piperidyl]-N-
15
               phenylpropanamide), its optical isomers, salts, and
16
               salts of isomers;
17
        [\frac{(10)}{(14)}] (14) Betacetylmethadol;
18
        \left[\frac{(11)}{(11)}\right] (15) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-
19
               phenethyl) -4-piperidinyl] -N-phenylpropanamide);
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1
        [\frac{(12)}{(16)}] (16) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-
2
               2-phenethyl)-3-methyl-4-piperidinyl]-N-
3
               phenylpropanamide);
4
        (17) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-
5
               2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide);
6
        [\frac{(13)}{(18)}] (18) Betameprodine;
7
        [(14)] (19) Betamethadol;
8
        (20) Beta-methyl fentanyl (N-phenyl-N-(1-(2-
9
               phenylpropyl)piperidin-4-yl)propionamide) (Other name:
10
               [beta] - methyl fentanyl);
11
        (21) Beta'-phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-
12
               N, 3-diphenylpropanamide) (Other names: [beta]'-phenyl
13
               fentanyl; 3-phenylpropanoyl fentanyl);
14
        [\frac{(15)}{(15)}] (22) Betaprodine;
15
        (23) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
16
               phenylbutyramide);
17
        [<del>(16)</del>] (24) Clonitazene;
18
        (25) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
19
               phenylcyclopropanecarboxamide);
20
        [\frac{17}{17}] (26) Dextromoramide;
21
        [\frac{(18)}{(18)}] (27) Diampromide;
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1
        [\frac{(19)}{(19)}] (28) Diethylthiambutene;
2
        \left[\frac{(20)}{(29)}\right] (29) Difenoxin;
 3
        [\frac{(21)}{(21)}] (30) Dimenoxadol;
 4
        [\frac{(22)}{(31)}] (31) Dimepheptanol;
 5
        [<del>(23)</del>] (32) Dimethylthiambutene;
 6
        [(24)] (33) Dioxaphetyl butyrate;
 7
        \left[\frac{(25)}{(25)}\right] (34) Dipipanone;
8
        [\frac{(26)}{}] (35) Ethylmethylthiambutene;
9
        \left[\frac{(27)}{}\right] (36) Etonitazene;
10
        [\frac{(28)}{(28)}] (37) Etoxeridine;
11
         (38) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-
12
                yl) (phenyl) carbamate);
13
         (39) 4-fluoroisobutyryl fentanyl [N-(4-fluorophenyl)-N-(1-
14
                phenethylpiperidin-4-yl)isobutyramide] (Other name:
15
                 para-fluoroisobutyryl fentanyl);
16
         (40) 2'-fluoro ortho-fluorofentanyl (N-(1-(2-
17
                 fluorophenethyl)piperidin-4-yl)-N-(2-
18
                 fluorophenyl)propionamide) (Other name: 2'-fluoro 2-
19
                 fluorofentanyl);
20
         (41) Furanyl fentanyl (N-(1-phenethylpiperidin-4-y1)-N-
21
                 phenylfuran-2-carboxamide);
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1
        [\frac{(29)}{(42)}] (42) Furethidine;
2
        [(30)] (43) Hydroxypethidine;
 3
        [-(31)] (44) Ketobemidone;
 4
        [\frac{(32)}{(32)}] (45) Levomoramide;
 5
        [<del>(33)</del>] (46) Levophenacylmorphan;
 6
         (47) Methoxyacetyl fentanyl (2-methoxy-N-(1-
 7
                phenethylpiperidin-4-yl) -N-phenylacetamide);
8
         (48) 4'-methyl acetyl fentanyl (N-(1-(4-
9
                methylphenethyl)piperidin-4-yl)-N-phenylacetamide);
10
        [\frac{34}{34}] (49) [\frac{3-\text{Methylfentanyl}}{3-\text{methylfentanyl}}] 3-methylfentanyl (N-[3-
11
                 methyl-1-(2-phenylethyl)-4-piperidyl]-N-
12
                phenylpropanamide);
13
        [\frac{35}{3}] (50) 3-methylthiofentanyl (N-[3-methyl-1-(2-
14
                 thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
15
        [(36)] (51) Morpheridine;
16
        [<del>(37)</del>] (52) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
17
        [(38)] (53) Noracymethadol;
18
        [\frac{(39)}{(39)}] (54) Norlevorphanol;
19
        [\frac{40}{1}] (55) Normethadone;
20
        [\frac{(41)}{(56)}] (56) Norpipanone;
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1	<u>(57)</u>	Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-
2		<pre>phenethylpiperidin-4-yl)acetamide];</pre>
3	(58)	Ortho-fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-
4		<pre>phenethylpiperidin-4-yl)acrylamide);</pre>
5	<u>(59)</u>	Ortho-fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-
6		phenethylpiperidin-4-yl)butyramide) (Other name: 2-
7		<pre>fluorobutyryl fentanyl);</pre>
8	(60)	Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-
9		phenethylpiperidin-4-yl)propionamide) (Other name: 2-
10		<pre>fluorofentanyl);</pre>
11	(61)	Ortho-fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-
12		(1-phenethylpiperidin-4-yl)isobutyramide);
13	(62)	Ortho-methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-
14		phenethylpiperidin-4-yl)acetamide) (Other name: 2-
15		<pre>methyl acetylfentanyl);</pre>
16	(63)	Ortho-methyl methoxyacetyl fentanyl (2-methoxy-N-(2-
17		methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide)
18		(Other name: 2-methyl methoxyacetyl fentanyl);
19	(64)	Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-
20		<pre>phenethylpiperidin-4-yl)butyramide);</pre>

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1
        [\frac{(42)}{(42)}] (65) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-
2
               phenethyl) -4-piperidinyl]propanamide;
3
        (66) Para-fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-
4
               phenethylpiperidin-4-yl) furan-2-carboxamide);
5
        (67) Para-methylfentanyl (N-(4-methylphenyl)-N-(1-
6
               phenethylpiperidin-4-yl)propionamide) (Other name: 4-
7
               methylfentanyl);
8
        [(43)] (68) PEPAP (1-(-2-phenethyl)-4-phenyl-4-
9
               acetoxypiperidine(+));
10
        [(44)] (69) Phenadoxone;
11
        [(45)] (70) Phenampromide;
       [(46)] (71) Phenomorphan;
12
13
        [\frac{(47)}{}] (72) Phenoperidine;
14
        (73) Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
15
               phenylbenzamide) (Other name: benzoyl fentanyl);
16
        [(48)] (74) Piritramide;
17
       [\frac{(49)}{(75)}] (75) Proheptazine;
       [(50)] (76) Properidine;
18
19
        \left[\frac{(51)}{(77)}\right] (77) Propiram;
20
        [\frac{(52)}{(78)}] Racemoramide;
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1
        (79)
               Thenylfentanyl (N-[1-(2-thienyl)methyl-4-piperidyl]-N-
2
               phenylpropanamide), its optical isomers, salts, and
3
               salts of isomers;
4
        \lceil \frac{(53)}{(53)} \rceil (80) Thiofentanyl (N-phenyl-N-\lceil 1-(2-\text{thienyl}) \text{ ethyl}-4-
5
               piperidinyl]-propanamide);
6
        (81) Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
7
               phenylthiophene-2-carboxamide) (Other names: 2-
8
               thiofuranyl fentanyl; thiophene fentanyl);
9
        [+54] (82) Tilidine;
10
        \lceil \frac{(55)}{(55)} \rceil (83) Trimeperidine; and
11
        [(56) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
12
               (benzylfentanyl), its optical isomers, salts, and
13
               salts of isomers;
14
        (57) N-[1-(2-thienyl)methyl-4-piperidyl]-N-
15
               phenylpropanamide (thenylfentanyl), its optical
16
               isomers, salts, and salts of isomers;
17
        (58) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide,
18
               (acetyl fentanyl), its optical, positional, and
19
               geometric isomers, salts, and salts of isomers;
20
        (59) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)
21
               eyclohexylmethyl]benzamide), its isomers, esters,
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ethers, salts, and salts of isomers, esters, and
1
2
              ethers;
       (60) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, its
3
              isomers, esters, ethers, salts, and salts of isomers,
4
              esters, and ethers (Other names: Butyryl fentanyl);
5
        (61) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-
6
              vll-N-phenylpropionamide, its isomers, esters, ethers,
7
              salts and salts of isomers, esters, and ethers (Other
8
9
              names: beta-hydroxythiofentanyl);
        (62) N-(1-phenethylpiperidin-4-y1)-N-phenylfuran-2-
10
              carboxamide, its isomers, esters, ethers, salts, and
11
              salts of isomers, esters, and ethers (Other names:
12
13
              Furanyl fentanyl);
        (63)] (84) U-47700 (3,4-dichloro-N-[2-
14
              (dimethylamino) cyclohexyl] -N-methylbenzamide[, its
15
              isomers, esters, ethers, salts, and salts of isomers,
16
17
              esters, and ethers (Other names: U-47700);
18
        (64) 4-fluoroisobutyryl fentanyl or para-fluoroisobutyryl
19
              fentanyl [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-
20
              4-yl) isobutyramide];
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1	(65)	Acryl fentanyl or acryloylfentanyl [N-(1-
2		<pre>phenethylpiperidin-4-yl) -N-phenylacrylamide];</pre>
3	(66)	Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-
4		<pre>phenethylpiperidin-4-yl) acetamide];</pre>
5	(67)	Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
6		phenylcyclopropanecarboxamide;
7	(68)	Methoxyacetyl fentanyl (2-methoxy-N-(1-
8		<pre>phenethylpiperidin-4-yl)-N-phenylacetamide);</pre>
9	(69)	Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-
10		phenethylpiperidin-4-yl)propionamide) (Other name: 2-
11		fluorofentanyl); and
12	(70)	Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-
13		<pre>phenethylpiperidin-4-yl)butyramide])."</pre>
14	2.	By amending subsections (f) and (g) to read:
15	"(f)	Stimulants. Unless specifically excepted or unless
16	listed in	another schedule, any material, compound, mixture, or
17	preparati	on [which] that contains any quantity of the following
18	substance	s having a stimulant effect on the central nervous
19	system, i	ncluding its salts, isomers, and salts of isomers:
20	(1)	Aminorex;
21	(2)	Cathinone;

1	(3)	4,4'-dimethylaminorex (common name: 4,4'-DMAR);
2	[-(3) -]	(4) Fenethylline;
3	[(4)]	(5) Methcathinone;
4	[-(5)	N-cthylamphetamine;
5	(6)	4-methylaminorex;
6	<u>(7)</u>	N-ethylamphetamine;
7	[-(7)-]	(8) N,N-dimethylamphetamine; [and]
8	[(8)]	(9) Substituted cathinones, any compound, except
9		bupropion or compounds listed under a different
10		schedule, structurally derived from 2-aminopropan-1-
11		one by substitution at the 1-position with either
12		phenyl, naphthyl, or thiophene ring systems, whether
13		or not the compound is further modified in any of the
14		following ways:
15		(A) By substitution in the ring system to any extent
16		with alkyl, alkylenedioxy, alkoxy, haloalkyl,
17		hydroxyl, or halide substituents, whether or not
18		further substituted in the ring system by one or
19		more other univalent substituents;
20		(B) By substitution at the 3-position with an acyclic
21		alkyl substituent; or

1	(C) By substitution at the 2-amino nitrogen atom with
2	alkyl, dialkyl, benzyl, or methoxybenzyl groups,
3	or by inclusion of the 2-amino nitrogen atom in a
4	cyclic structure.
5	Some other trade names: Mephedrone (2-methylamino-1-p-
6	tolylpropan-1-one), also known as 4-
7	methylmethcathinone (4-MMC), methylephedrone or MMCAT;
8	Methylenedioxypyrovalerone (MDPV, MDPK); methylone or
9	3,4-methylenedioxymethcathinone; and 1-
10	(benzo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one,
11	monohydrochloride, also known as Ethylone, bk-MDEA
12	hydrochloride, MDEC; 3,4-Methylenedioxy-N-
13	ethylcathinone; bk-Methylenedioxyethylamphetamine, 4-
14	methyl-N-ethylcathinone (4-MEC); 4-methyl-alpha-
15	pyrrolidinopropiophenone (4-MePPP); alpha-
16	<pre>pyrrolidinopentiophenone ([alpha]-PVP); 1-(1,3-</pre>
17	benzodioxol-5-yl)-2-(methylamino)butan-1-one
18	(butylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-
19	1-one (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-
20	(methylamino)pentan-1-one (pentylone, bk-MBDP); 4-
21	fluoro-N-methylcathinone (4-FMC, flephedrone); 3-

1		fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-
2		2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-
3		pyrrolidinobutiophenone ([alpha]-PBP) and their
4		optical, positional, and geometric isomers, salts and
5		salts of isomers, whenever the existence of such
6		salts, isomers, and salts of isomers is possible $[\cdot]$:
7		and
8	(10)	1-(4-methoxyphenyl)-N-methylpropan-2-amine (Other
9		names: para-methoxymethamphetamine; PMMA).
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
18		resinous extractives of Cannabis, sp. or synthetic
19		substances, derivatives, and their isomers with
20		similar chemical structure and pharmacological
21		activity to those substances contained in the plant,

1		such as the following: Delta I cls or trans
2		tetrahydrocannabinol, and their optical isomers; Delta
3		6 cis or trans tetrahydrocannabinol, and their optical
4		isomers; and Delta 3,4 cis or trans-
5		tetrahydrocannabinol, and its optical isomers (since
6		nomenclature of these substances is not
7		internationally standardized, compounds of these
8		structures, regardless of numerical designation of
9		atomic positions, are covered);
10	(2)	Naphthoylindoles; meaning any compound containing a 3-
11		(1-naphthoyl) indole structure with substitution at the
12		nitrogen atom of the indole ring by a alkyl,
13		haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
14		1-(N-methyl-2-piperidinyl) methyl or 2-(4-
15		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	(3)	Naphthylmethylindoles; meaning any compound containing
20		a 1H-indol-3-yl-(1-naphthyl) methane structure with
21		substitution at the nitrogen atom of the indole ring

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

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

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1
              1-(N-methyl-2-piperidinyl) methyl, or 2-(4-
2
              morpholinyl) ethyl group whether or not further
3
              substituted in the indole ring to any extent and
4
              whether or not substituted in the phenyl ring to any
5
              extent;
6
         (9)
              [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
7
              pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-
8
              naphthalenylmethanone (another trade name is WIN
9
              55,212-2);
10
        (10)
              (6a, 10a) - 9 - (hydroxymethyl) - 6, 6 - dimethyl - 3 - (2 - 6)
11
              methyloctan-2-yl)-6a,7,10,10a-
12
              tetrahydrobenzo[c]chromen-1-ol (Other trade names are:
13
              HU-210/HU-211);
14
        (11)
              Tetramethylcyclopropanoylindoles; meaning any compound
15
              containing a 3-tetramethylcyclopropanoylindole
16
              structure with substitution at the nitrogen atom of
17
              the indole ring by an alkyl, haloalkyl, cyanoalkyl,
18
              alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
19
              methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl,
20
              1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
21
              morpholinyl) methyl, or tetrahydropyranylmethyl group,
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1
               whether or not further substituted in the indole ring
2
               to any extent and whether or not substituted in the
3
               tetramethylcyclopropyl ring to any extent;
4
        (12)
              N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
5
               its optical, positional, and geometric isomers, salts,
 6
               and salts of isomers (Other names: APINACA, AKB48);
7
        (13)
              Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
8
               optical, positional, and geometric isomers, salts, and
9
               salts of isomers (Other names: PB-22; QUPIC);
10
         (14)
               Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-
11
               carboxylate, its optical, positional, and geometric
12
               isomers, salts, and salts of isomers (Other names: 5-
13
               fluoro-PB-22; 5F-PB-22);
14
        (15)
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-amino-3-methyl-1-oxobutan-2-yl)
15
               fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
16
               positional, and geometric isomers, salts, and salts of
17
               isomers (Other names: AB-FUBINACA);
18
        (16)
              N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
19
               indazole-3-carboxamide, its optical, positional, and
20
               geometric isomers, salts, and salts of isomers (Other
21
               names: ADB-PINACA);
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```
1
             N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
        (17)
2
              (cyclohexylmethyl) -1H-indazole-3-carboxamide, its
              optical, positional, and geometric isomers, salts, and
3
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)-1H-indazole-3-carbonyl)-L-
12
              valinate, and geometric isomers, salts, and salts of
13
              isomers (Other names: FUB-AMB, Methyl 2-(1-(4-
14
              fluorobenzyl) -1H-indazole-3-carboxamido) -3-
15
              methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);
16
        (21)
             (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
              carboxamido) -3-methylbutanoate, and geometric isomers,
17
18
              salts, and salts of isomers (Other names: 5-fluoro-
19
              AMB, 5-fluoro-AMP);
20
        (22)
              N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-
21
              indazole-3-carboxamide, and geometric isomers, salts,
```

1		and salts of isomers (Other names: AKB48 N-(5-
2		fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
3		analog, 5F-APINACA);
4	(23)	N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
5		geometric isomers, salts, and salts of isomers (Other
6		names: STS-135, 5F-APICA; 5-fluoro-APICA);
7	(24)	Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
8		carboxylate, and geometric isomers, salts, and salts
9		of isomers (Other names: NM2201[+]; CBL2201);
10	(25)	N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
11		(cyclohexylmethyl)-1H-indazole-3-carboxamide, and
12		geometric isomers, salts, and salts of isomers (Other
13		names: MAB-CHMINACA and ADB-CHMINACA);
14	(26)	Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
15		carboxamido]-3,3-dimethylbutanoate (Other names: 5F-
16		ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical,
17		positional, and geometric isomers, salts, and salts of
18		isomers; [and]
19	(27)	[1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)indazole-3-
20		carboxamide (CUMYL-4CN-BINACA), 1-(4-cyanobutyl)-N-
21		(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its

1		optical, positional, and geometric isomers, salts, and
2		salts of isomers[; also known as] (Other names: SGT-
3		78[, 4-CN-CUMYL-BINACA]; <u>4-CN-CUMYL BINACA; 4-CN-</u>
4		CUMYL-BUTINACA; CUMYL-CB-PINACA; CUMYL-CYBINACA; [4-
5		eyano CUMYL-BUTINACA.] 4-cyano-CUMYL-BUTINACA; CUMYL-
6		4CN-BINACA);
7	(28)	N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
8		fluoropentyl)-1H-indazole-3-carboxamide (Other name:
9		5F-AB-PINACA);
10	(29)	Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
11		carboxamido)-3-methylbutanoate (Other names: MMB-
12		CHMICA; AMB-CHMICA);
13	(30)	1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
14		<pre>pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-</pre>
15		CUMYL-P7AICA); and
16	(31)	Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
17		3-carboxamido)butanoate (MDMB-4en-PINACA)."
18	SECT	ION 2. Section 329-16, Hawaii Revised Statutes, is
19	amended b	y amending subsection (c) to read as follows:
20	"(c)	Any of the following opiates, including their
21	isomers.	esters, ethers, salts, and salts of isomers, whenever

```
1
    the existence of these isomers, esters, ethers, and salts is
2
    possible within the specific chemical designation:
3
         (1)
              Alfentanil;
4
         (2)
              Alphaprodine;
5
         (3) Anileridine;
 6
         (4) Bezitramide;
7
         (5)
              Bulk Dextropropoxyphene (nondosage form);
8
              Carfentanil;
         (6)
9
         (7)
              Dihydrocodeine;
10
         (8)
              Diphenoxylate;
11
         (9)
              Fentanyl;
12
        (10)
              Isomethadone;
13
        (11)
              Levo-alphacetylmethadol (LAAM);
14
        (12)
              Levomethorphan;
15
        (13)
              Levorphanol;
16
        (14)
              Metazocine;
17
        (15)
              Methadone;
18
        (16)
              Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-
19
              diphenyl butane;
20
              Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-
        (17)
21
              diphenyl-propane-carboxylic acid;
```

```
1
           (18) Oliceridine, including the free base form, and its
  2
                  salts, to include the fumarate salt, by definition;
  3
          [<del>(18)</del>] (19) Pethidine (Meperidine);
  4
          [<del>(19)</del>] (20) Pethidine-Intermediate-A, 4-cyano-1-methyl-4-
  5
                  phenylpiperidine;
  6
          [<del>(20)</del>] (21) Pethidine-Intermediate-B, ethyl-4-
  7
                  phenylpiperidine-4-carboxylate;
  8
          [<del>(21)</del>] (22) Pethidine-Intermediate-C, 1-methyl-4-
  9
                  phenylpiperidine-4-carboxylic acid;
 10
          \left[\frac{(22)}{(23)}\right] (23) Phenazocine;
 11
          \left[\frac{(23)}{(24)}\right] (24) Piminodine;
 12
          [\frac{(24)}{}] (25) Racemethorphan;
 13
          [\frac{(25)}{}] (26) Racemorphan;
 14
          [\frac{(26)}{}] (27) Remifentanil;
15
         \left[\frac{(27)}{(28)}\right] (28) Sufentanil;
          [\frac{(28)}{(29)}] (29) Tapentadol; and
 16
          [\frac{(29)}{(30)}] (30) Thiafentanil."
 17
            SECTION 3. Section 329-20, Hawaii Revised Statutes, is
 18
 19
      amended as follows:
```

1. By amending subsection (b) to read:

20

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1
          "(b) Depressants. Any material, compound, mixture, or
 2
    preparation which contains any quantity of the following
 3
     substances, including its salts, isomers, esters, ethers, and
 4
    salts of isomers, whenever the existence of these isomers,
 5
    esters, ethers, and salts is possible within the specific
 6
    chemical designation, that has a degree of danger or probable
 7
    danger associated with a depressant effect on the central
 8
    nervous system:
 9
           (1)
               Alprazolam;
10
           (2) Barbital;
11
          (3) Brexanolone;
12
         [-(3)] (4) Bromazepam;
13
         [(4)] (5) Butorphanol;
14
         [-(5)] (6) Camazepam;
15
         \left[\frac{(6)}{(7)}\right] (7) Carisoprodol;
16
         [\frac{(7)}{1}] (8) Chloral betaine;
17
         [<del>(8)</del>] (9) Chloral hydrate;
18
         [\frac{(9)}{(10)}] (10) Chlordiazepoxide;
19
        \left[\frac{(10)}{(11)}\right] (11) Clobazam;
20
        [<del>(11)</del>] <u>(12)</u> Clonazepam;
21
        [\frac{(12)}{(13)}] (13) Clorazepate;
```

```
1
         [<del>(13)</del>] (14) Clotiazepam;
 2
         [\frac{(14)}{(15)}] (15) Cloxazolam;
 3
         [\frac{(15)}{(16)}] (16) Delorazepam;
 4
         (16) Dichloralphenazone (Midrin);
 5
         (17) Diazepam;
 6
          (18) Dichloralphenazone (Midrin);
 7
         \left[\frac{(18)}{(19)}\right] (19) Estazolam;
         [\frac{(19)}{(20)}] Ethchlorvynol;
 8
 9
         [\frac{(20)}{(21)}] (21) Ethinamate;
10
         [\frac{(21)}{(21)}] (22) Ethyl loflazepate;
11
         \left[\frac{(22)}{(23)}\right] (23) Fludiazepam;
12
         [\frac{(23)}{(24)}] (24) Flunitrazepam;
13
         [-(24)] (25) Flurazepam;
14
         [\frac{(25)}{(26)}] (26) Fospropofol (Lusedra);
15
         \left[\frac{(26)}{(27)}\right] (27) Halazepam;
16
         \left[\frac{(27)}{(28)}\right] (28) Haloxazolam;
17
         [\frac{(28)}{(29)}] (29) Ketazolam;
18
          (30) Lemborexant ((1R,2S)-2-[(2,4-dimethylpyrimidin-5-
19
                   yl)oxymethyl]-2-(3-fluorophenyl)-N-(5-fluoropyridin-2-
20
                   yl)cyclopropane-1-carboxamide);
21
         [<del>(29)</del>] (31) Loprazolam;
```

```
1
         [(30)] (32) Lorazepam;
2
         [(31)] (33) Lormetazepam;
3
         [\frac{(32)}{(34)}] Mebutamate;
 4
         [<del>(33)</del>] (35) Medazepam;
 5
         [<del>(34)</del>] (36) Meprobamate;
         [<del>(35)</del>] <u>(37)</u> Methohexital;
 6
7
         [<del>(36)</del>] (38) Methylphenobarbital (mephorbarbital);
8
         [<del>(37)</del>] <u>(</u>39) Midazolam;
9
         [\frac{(38)}{(40)}] (40) Nimetazepam;
10
         [\frac{(39)}{(41)}] Nitrazepam;
         [\frac{(40)}{(42)}] Nordiazepam;
11
12
         \left[\frac{(41)}{(41)}\right] (43) Oxazepam;
13
         [\frac{(42)}{(42)}] (44) Oxazolam;
         [<del>(43)</del>] (45) Paraldehyde;
14
         [<del>(44)</del>] (46) Petrichloral;
15
16
         [(45)] (47) Phenobarbital;
17
         [-(46)] (48) Pinazepam;
18
         [-(47)] (49) Prazepam;
19
         [-(48)] (50) Quazepam;
20
         (51) Remimazolam;
21
         [(49)] (52) Suvorexant;
```

```
1
        \left[\frac{(50)}{(53)}\right] (53) Temazepam;
2
        [(51)] (54) Tetrazepam;
3
        [<del>(52)</del>] <u>(55)</u> Triazolam;
4
        \left[\frac{(53)}{(56)}\right] (56) Zaleplon;
5
        [(54)] (57) Zolpidem; and
6
        [(55)] (58) Zopiclone (Lunesta) [; and
7
        (56) Brexanolone]."
8
          2. By amending subsection (d) to read:
          "(d) Stimulants. Unless listed in another schedule, any
9
10
    material, compound, mixture, or preparation which contains any
11
    quantity of the following substances having a stimulant effect
12
    on the central nervous system, including its salts, isomers, and
13
    salts of such isomers whenever the existence of such salts,
14
    isomers, and salts of isomers is possible within the specific
    chemical designation:
15
16
          (1)
                Cathine ((+)-norpseudoephedrine);
17
          (2)
               Diethylpropion;
18
          (3)
               Fencamfamin;
19
          (4)
               Fenproporex;
20
          (5) Lorcaserin;
21
         \left[\frac{(5)}{(6)}\right] (6) Mazindol;
```

```
1
         [\frac{(6)}{(7)}] (7) Mefenorex;
2
         \left[\frac{7}{1}\right] (8) Modafinil;
3
         [<del>(8)</del> Phentermine;
4
              Pemoline (including organometallic complexes and
          (9)
5
               chelates thereof);
6
        (10) Phentermine;
7
        [<del>(10)</del>] (11) Pipradrol;
8
        (12) Serdexmethylphenidate;
9
        [<del>(11)</del>] (13) Sibutramine;
10
        (14) Solriamfetol; and
11
        [\frac{(12)}{(15)}] (15) SPA (1-dimethylamino-1,2-diphenylethane,
12
               lefetamine) [+
13
        (13) Lorcaserin; and
14
        (14) Solriamfetol]."
15
          SECTION 4. Section 329-22, Hawaii Revised Statutes, is
16
    amended by amending subsection (d) to read as follows:
17
          "(d) Depressants. Unless specifically exempted or
    excluded or unless listed in another schedule, any material,
18
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) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-
4
              yl]butanamide) (Other names: BRV; UCB-34714; Briviact)
5
              and its salts;
        [\frac{1}{1}] (2) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-
6
7
              propionamide], (Vimpat);
8
         (3) Lasmiditan (2,4,6-trifluoro-N-(6-(1-methylpiperidine-
9
              4-carbonyl)pyridine-2-yl-benzamide); and
10
        [+(2)] (4) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic
11
              acid][; and
12
         (3) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-
13
              yl]butanamide) (Other names: BRV; UCB-34714; Briviact)
14
              and its salts]."
15
         SECTION 5. Statutory material to be repealed is bracketed
16
    and stricken. New statutory material is underscored.
17
         SECTION 6. This Act shall take effect on January 1, 2050.
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Report Title:

Uniform Controlled Substances Act

Description:

Amends the schedules for controlled substances under the Uniform Controlled Substances Act for purposes of conforming with updates in federal law. Effective 1/1/2050. (HD2)

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