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 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 Acetyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-(3) 13 phenylacetamide), its optical, positional, and 14 geometric isomers, salts, and salts of isomers; 15 (4) Acryl fentanyl [N-(1-phenethylpiperidin-4-yl)-N-16 phenylacrylamide] (Other name: acryloylfentanyl); 17 (5) AH-7921 (3,4-dichloro-N-[(1-18 dimethylamino)cyclohexylmethyl]benzamide);

```
1
         [<del>(3)</del>] (6) Allylprodine;
          [\frac{4}{1}] (7) Alphacetylmethadol (except levo-
 2
 3
                alphacetylmethadol, levomethadyl acetate, or LAAM);
          [<del>(5)</del>] (8) Alphameprodine;
 4
5
         [\frac{(6)}{(9)}] (9) Alphamethadol;
          [<del>(7)</del>] (10) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-
 6
 7
                phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-
8
                2-phenylethyl) -4-(N-propanilido)piperidine);
9
          [\frac{(8)}{(11)}] (11) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-
10
                thienyl) ethyl-4-piperidinyl] -N-phenylpropanamide);
         [\frac{(9)}{}] (12) Benzethidine;
11
12
                Benzylfentanyl (N-[1-benzyl-4-piperidyl]-N-
         (13)
                phenylpropanamide), its optical isomers, salts, and
13
14
                salts of isomers;
15
        [<del>(10)</del>] (14) Betacetylmethadol;
16
        \left[\frac{(11)}{(15)}\right] (15) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-
17
                phenethyl) -4-piperidinyl] -N-phenylpropanamide);
18
        \left[\frac{(12)}{(16)}\right] (16) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-
19
                2-phenethyl)-3-methyl-4-piperidinyl]-N-
20
                phenylpropanamide);
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1
          (17) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-
 2
                 2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide);
 3
         [\frac{(13)}{(18)}] (18) Betameprodine;
 4
         \left[\frac{14}{14}\right] (19) Betamethadol;
 5
         (20) Beta-methyl fentanyl (N-phenyl-N-(1-(2-
 6
                 phenylpropyl)piperidin-4-yl)propionamide) (Other name:
 7
                 [beta] -methyl fentanyl);
 8
         (21) Beta'-phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-
 9
                 N, 3-diphenylpropanamide) (Other names: [beta]'-phenyl
10
                 fentanyl; 3-phenylpropanoyl fentanyl);
11
         \left[\frac{(15)}{(15)}\right] (22) Betaprodine;
12
         (23) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
13
                 phenylbutyramide);
14
         [<del>(16)</del>] (24) Clonitazene;
15
         (25) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
16
                 phenylcyclopropanecarboxamide);
17
        [\frac{(17)}{(17)}] (26) Dextromoramide;
18
        [\frac{(18)}{(27)}] (27) Diampromide;
19
        [<del>(19)</del>] (28) Diethylthiambutene;
20
        \left[\frac{(20)}{(20)}\right] (29) Difenoxin;
21
        \left[\frac{(21)}{(21)}\right] (30) Dimenoxadol;
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1
         [\frac{(22)}{(31)}] (31) Dimepheptanol;
 2
         [<del>(23)</del>] (32) Dimethylthiambutene;
 3
         [\frac{(24)}{(24)}] (33) Dioxaphetyl butyrate;
 4
         \left[\frac{(25)}{(25)}\right] (34) Dipipanone;
 5
         [\frac{(26)}{(26)}] (35) Ethylmethylthiambutene;
 6
         \left[\frac{(27)}{(27)}\right] (36) Etonitazene;
 7
         \left[\frac{(28)}{(28)}\right] (37) Etoxeridine;
 8
          (38) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-
 9
                 yl) (phenyl) carbamate);
10
          (39) 4-fluoroisobutyryl fentanyl [N-(4-fluorophenyl)-N-(1-
11
                 phenethylpiperidin-4-yl)isobutyramide] (Other name:
12
                 para-fluoroisobutyryl fentanyl);
13
          (40)
                 2'-fluoro ortho-fluorofentanyl (N-(1-(2-
14
                 fluorophenethyl)piperidin-4-yl)-N-(2-
15
                 fluorophenyl)propionamide) (Other name: 2'-fluoro 2-
16
                 fluorofentanyl);
17
          (41) Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
18
                 phenylfuran-2-carboxamide);
19
         \left[\frac{(29)}{(42)}\right] (42) Furethidine;
20
         [<del>(30)</del>] (43) Hydroxypethidine;
21
         [\frac{(31)}{(31)}] (44) Ketobemidone;
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1
         [\frac{(32)}{(32)}] (45) Levomoramide;
 2
         [<del>(33)</del>] (46) Levophenacylmorphan;
 3
                 Methoxyacetyl fentanyl (2-methoxy-N-(1-
          (47)
 4
                 phenethylpiperidin-4-yl) -N-phenylacetamide);
 5
          (48) 4'-methyl acetyl fentanyl (N-(1-(4-
 6
                 methylphenethyl)piperidin-4-yl)-N-phenylacetamide);
 7
         [\frac{(34)}{(34)}] (49) [\frac{3-\text{Methylfentanyl}}{(3-\text{Methylfentanyl}}] 3-methylfentanyl (N-[3-
 8
                 methyl-1-(2-phenylethyl)-4-piperidyl]-N-
 9
                 phenylpropanamide);
         \left[\frac{35}{3}\right] (50) 3-methylthiofentanyl (N-[3-methyl-1-(2-
10
11
                 thienyl) ethyl-4-piperidinyl] -N-phenylpropanamide);
12
         [\frac{(36)}{(51)}] (51) Morpheridine;
13
         [<del>(37)</del>] (52) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
14
         [\frac{(38)}{(53)}] (53) Noracymethadol;
15
         [\frac{(39)}{(39)}] (54) Norlevorphanol;
16
         [\frac{(40)}{}] (55) Normethadone;
17
         [<del>(41)</del>] (56) Norpipanone;
18
         (57)
                 Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-
19
                 phenethylpiperidin-4-yl)acetamide];
20
                 Ortho-fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-
         (58)
21
                 phenethylpiperidin-4-yl)acrylamide);
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1	(59)	Ortho-fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-
2		phenethylpiperidin-4-yl)butyramide) (Other name: 2-
3		<pre>fluorobutyryl fentanyl);</pre>
4	(60)	Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-
5		phenethylpiperidin-4-yl)propionamide) (Other name: 2-
6		<pre>fluorofentanyl);</pre>
7	(61)	Ortho-fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-
8		(1-phenethylpiperidin-4-yl) isobutyramide);
9	(62)	Ortho-methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-
10		phenethylpiperidin-4-yl)acetamide) (Other name: 2-
11		<pre>methyl acetylfentanyl);</pre>
12	(63)	Ortho-methyl methoxyacetyl fentanyl (2-methoxy-N-(2-
13		methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide)
14		(Other name: 2-methyl methoxyacetyl fentanyl);
15	(64)	Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-
16		<pre>phenethylpiperidin-4-yl)butyramide);</pre>
17	[(42)]	(65) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-
18		<pre>phenethyl) -4-piperidinyl]propanamide;</pre>
19	<u>(66)</u>	Para-fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-
20		<pre>phenethylpiperidin-4-yl)furan-2-carboxamide);</pre>

```
1
          (67) Para-methylfentanyl (N-(4-methylphenyl)-N-(1-
 2
                 phenethylpiperidin-4-yl)propionamide) (Other name: 4-
 3
                 methylfentanyl);
 4
         [\frac{(43)}{(43)}] (68) PEPAP (1-(-2-phenethyl)-4-phenyl-4-
 5
                 acetoxypiperidine[+]);
 6
         \left[\frac{44}{44}\right] (69) Phenadoxone;
 7
         [\frac{45}{1}] (70) Phenampromide;
         [-(46)-] (71) Phenomorphan;
 8
 9
         [\frac{47}{1}] (72) Phenoperidine;
10
         (73) Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
11
                 phenylbenzamide) (Other name: benzoyl fentanyl);
12
         [<del>(48)</del>] (74) Piritramide;
13
         [\frac{(49)}{(75)}] (75) Proheptazine;
14
         [<del>(50)</del>] (76) Properidine;
15
        \left[\frac{(51)}{}\right] (77) Propiram;
16
         \left[\frac{(52)}{(78)}\right] (78) Racemoramide;
17
                 Thenylfentanyl (N-[1-(2-thienyl)methyl-4-piperidyl]-N-
         (79)
18
                 phenylpropanamide), its optical isomers, salts, and
19
                 salts of isomers;
20
        [\frac{(53)}{(53)}] (80) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-
21
                 piperidinyl] -propanamide);
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Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
1
        (81)
               phenylthiophene-2-carboxamide) (Other names: 2-
2
               thiofuranyl fentanyl; thiophene fentanyl);
3
 4
       [<del>(54)</del>] (82) Tilidine;
5
       [\frac{(55)}{(55)}] (83) Trimeperidine; and
       [(56) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
6
 7
               (benzylfentanyl), its optical isomers, salts, and
8
               salts of isomers;
9
        (57) N-[1-(2-thienyl)methyl-4-piperidyl]-N-
               phenylpropanamide (thenylfentanyl), its optical
10
11
               isomers, salts, and salts of isomers;
              N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide,
12
        <del>(58)</del>
               (acetyl fentanyl), its optical, positional, and
13
14
               geometric isomers, salts, and salts of isomers;
15
        (59) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)
16
               cyclohexylmethyl]benzamide), its isomers, esters,
17
               ethers, salts, and salts of isomers, esters, and
18
              ethers;
19
        (60) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, its
20
               isomers, esters, ethers, salts, and salts of isomers,
21
               esters, and ethers (Other names: Butyryl fentanyl);
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1
        (61) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-
 2
               yl] N phenylpropionamide, its isomers, esters, ethers,
 3
               salts and salts of isomers, esters, and ethers (Other
 4
               names: beta-hydroxythiofentanyl);
 5
        <del>(62)</del>
              N-(1-phenethylpiperidin-4-y1)-N-phenylfuran-2-
 6
               carboxamide, its isomers, esters, ethers, salts, and
 7
               salts of isomers, esters, and others (Other names:
 8
              Furanyl fentanyl);
 9
        (63) (84) U-47700 (3,4-dichloro-N-[2-
10
               (dimethylamino) cyclohexyl] -N-methylbenzamide[, its
11
              isomers, esters, ethers, salts, and salts of isomers,
12
              esters, and ethers (Other names: U-47700);
13
        (64) 4-fluoroisobutyryl fentanyl or para-fluoroisobutyryl
14
              fentanyl [N-(4-fluorophenyl) N-(1-phenethylpiperidin-
15
              4-yl) isobutyramide];
16
        (65) Acryl fentanyl or acryloylfentanyl [N-(1-
17
              phenethylpiperidin-4-vl)-N-phenylacrylamidel;
18
        (66) Ocfentanil [N-(2-fluorophenyl) 2-methoxy-N-(1-
19
              phenethylpiperidin-4-yl)acetamide];
20
        (67) Cyclopropyl-fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
21
              phenylcyclopropanecarboxamide;
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```
1
        (68) Methoxyacetyl fentanyl (2-methoxy-N (1-
              phenethylpiperidin-4-yl) -N-phenylacetamide);
2
3
        (69) Ortho-fluorofentanyl (N-(2-fluorophenyl) N-(1-
              phenethylpiperidin-4-yl)propionamide) (Other name: 2-
4
              fluorofentanyl); and
5
        (70) Para-fluorobutyryl fentanyl (N (4-fluorophenyl)-N-(1-
6
7
              phenethylpiperidin-4-yl)butyramide])."
             By amending subsections (f) and (g) to read:
8
               Stimulants. Unless specifically excepted or unless
9
         "(f)
    listed in another schedule, any material, compound, mixture, or
10
11
    preparation [which] that contains any quantity of the following
12
    substances having a stimulant effect on the central nervous
    system, including its salts, isomers, and salts of isomers:
13
14
         (1)
              Aminorex;
15
         (2)
              Cathinone;
         (3) 4,4'-dimethylaminorex (common name: 4,4'-DMAR);
16
17
        \left[\frac{(3)}{(4)}\right] (4) Fenethylline;
18
        [\frac{4}{1}] (5) Methcathinone;
19
        [(5) N-ethylamphetamine;]
20
         (6) 4-methylaminorex;
21
         (7) N-ethylamphetamine;
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1	[(7)]	<u>(8)</u>	N,N-dimethylamphetamine; [and]
2	[-(8) -]	<u>(9)</u>	Substituted cathinones, any compound, except
3		bupr	opion or compounds listed under a different
4		sche	dule, structurally derived from 2-aminopropan-1-
5		one :	by substitution at the 1-position with either
6		phen	yl, naphthyl, or thiophene ring systems, whether
7		or n	ot the compound is further modified in any of the
8		foll	owing ways:
9		(A)	By substitution in the ring system to any extent
10			with alkyl, alkylenedioxy, alkoxy, haloalkyl,
11			hydroxyl, or halide substituents, whether or not
12			further substituted in the ring system by one or
13			more other univalent substituents;
14		(B)	By substitution at the 3-position with an acyclic
15			alkyl substituent; or
16		(C)	By substitution at the 2-amino nitrogen atom with
17			alkyl, dialkyl, benzyl, or methoxybenzyl groups,
18			or by inclusion of the 2-amino nitrogen atom in a
19			cyclic structure.
20		Some	other trade names: Mephedrone (2-methylamino-1-p-
21		toly	lpropan-1-one), also known as 4-

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1
              methylmethcathinone (4-MMC), methylephedrone or MMCAT;
              Methylenedioxypyrovalerone (MDPV, MDPK); methylone or
 2
 3
              3,4-methylenedioxymethcathinone; and 1-
               (benzo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one,
 4
 5
              monohydrochloride, also known as Ethylone, bk-MDEA
              hydrochloride, MDEC; 3,4-Methylenedioxy-N-
 6
 7
              ethylcathinone; bk-Methylenedioxyethylamphetamine, 4-
8
              methyl-N-ethylcathinone (4-MEC); 4-methyl-alpha-
9
              pyrrolidinopropiophenone (4-MePPP); alpha-
10
              pyrrolidinopentiophenone ([alpha]-PVP); 1-(1,3-
11
              benzodioxol-5-yl)-2-(methylamino)butan-1-one
12
              (butylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-
13
              1-one (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-
14
              (methylamino) pentan-1-one (pentylone, bk-MBDP); 4-
15
              fluoro-N-methylcathinone (4-FMC, flephedrone); 3-
16
              fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-
17
              2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-
18
              pyrrolidinobutiophenone ([alpha]-PBP) and their
19
              optical, positional, and geometric isomers, salts and
20
              salts of isomers, whenever the existence of such
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S.B. NO. 3141 S.D. 1 H.D. 2 C.D. 1

1		salts, isomers, and salts of isomers is possible $[-]$;
2		and
3	(10)	1-(4-methoxyphenyl)-N-methylpropan-2-amine (Other
4		names: para-methoxymethamphetamine; PMMA).
5	(g)	Any of the following cannabinoids, their salts,
6	isomers,	and salts of isomers, unless specifically excepted,
7	whenever	the existence of these salts, isomers, and salts of
8	isomers i	s possible within the specific chemical designation:
9	(1)	Tetrahydrocannabinols; meaning tetrahydrocannabinols
10		naturally contained in a plant of the genus Cannabis
11		(cannabis plant), as well as synthetic equivalents of
12		the substances contained in the plant, or in the
13		resinous extractives of Cannabis, sp. or synthetic
14		substances, derivatives, and their isomers with
15		similar chemical structure and pharmacological
16		activity to those substances contained in the plant,
17		such as the following: Delta 1 cis or trans
18		tetrahydrocannabinol, and their optical isomers; Delta
19		6 cis or trans tetrahydrocannabinol, and their optical
20		isomers; and Delta 3,4 cis or trans-
21		tetrahydrocannabinol, and its optical isomers (since

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

1

16

17

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19

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21

S.B. NO. S.D. 1 H.D. 2

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-morpholinyl)

whether or not substituted in the naphthyl ring to any

(6) Phenylacetylindoles; meaning any compound containing a 3-phenylacetylindole structure with substitution at the nitrogen atom of the indole ring by a alkyl,

in the naphthyl ring to any extent;

ethyl group whether or not further substituted in the

indene ring to any extent, whether or not substituted

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		indole ring to any extent, whether or not substituted
5		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,
16		haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
17		1-(N-methyl-2-piperidinyl) methyl, or 2-(4-
18		morpholinyl) ethyl group whether or not further
19		substituted in the indole ring to any extent and
20		whether or not substituted in the phenyl ring to any
21		extent;

S.B. NO. S.D. 1

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1
               [2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)
         (9)
 2
              pyrrolo[1,2,3-de]-1, 4-benzoxazin-6-yl]-1-
 3
              naphthalenylmethanone (another trade name is WIN
 4
               55,212-2);
 5
               (6a, 10a) -9-(hydroxymethyl) -6, 6-dimethyl-3-(2-
        (10)
              methyloctan-2-yl)-6a,7,10,10a-
 6
 7
               tetrahydrobenzo[c]chromen-1-ol (Other trade names are:
 8
              HU-210/HU-211);
              Tetramethylcyclopropanoylindoles; meaning any compound
 9
        (11)
10
              containing a 3-tetramethylcyclopropanoylindole
11
              structure with substitution at the nitrogen atom of
12
              the indole ring by an alkyl, haloalkyl, cyanoalkyl,
13
              alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
14
              methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl,
15
              1-(N-methyl-2-pyrrolidinyl) methyl, 1-(N-methyl-3-
16
              morpholinyl) methyl, or tetrahydropyranylmethyl group,
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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N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide,
1
        (12)
              its optical, positional, and geometric isomers, salts,
2
              and salts of isomers (Other names: APINACA, AKB48);
3
4
        (13)
              Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate, its
              optical, positional, and geometric isomers, salts, and
5
              salts of isomers (Other names: PB-22; QUPIC);
6
7
              Quinolin-8-yl 1-(5fluoropentyl)-1H-indole-3-
        (14)
8
              carboxylate, its optical, positional, and geometric
              isomers, salts, and salts of isomers (Other names: 5-
9
10
              fluoro-PB-22; 5F-PB-22);
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-
11
        (15)
              fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
12
              positional, and geometric isomers, salts, and salts of
13
14
              isomers (Other names: AB-FUBINACA);
15
        (16)
              N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
              indazole-3-carboxamide, its optical, positional, and
16
              geometric isomers, salts, and salts of isomers (Other
17
18
              names: ADB-PINACA);
19
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
        (17)
              (cyclohexylmethyl) -1H-indazole-3-carboxamide, its
20
```

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

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

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

```
the existence of these isomers, esters, ethers, and salts is
1
2
    possible within the specific chemical designation:
3
         (1)
              Alfentanil;
4
         (2)
              Alphaprodine;
         (3) Anileridine;
5
         (4) Bezitramide;
6
7
              Bulk Dextropropoxyphene (nondosage form);
         (5)
8
         (6)
              Carfentanil;
9
              Dihydrocodeine;
         (7)
10
              Diphenoxylate;
         (8)
11
         (9)
              Fentanyl;
12
        (10)
              Isomethadone;
13
              Levo-alphacetylmethadol (LAAM);
        (11)
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;
         [<del>(20)</del>] (21) Pethidine-Intermediate-B, ethyl-4-
 6
 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
         [<del>(23)</del>] (24) Piminodine;
12
         [\frac{(24)}{2}] (25) Racemethorphan;
13
         \left[\frac{(25)}{(25)}\right] (26) Racemorphan;
14
         [\frac{(26)}{}] (27) Remifertanil;
15
         \left[\frac{(27)}{(28)}\right] (28) Suffentianil;
16
         [\frac{(28)}{(29)}] (29) Tapentadol; and
17
         [<del>(29)</del>] (30) Thiafentanil."
18
           SECTION 3. Section 329-20, Hawaii Revised Statutes, is
19
     amended as follows:
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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
     substances, including its salts, isomers, esters, ethers, and
 3
 4
     salts of isomers, whenever the existence of these isomers,
 5
     esters, ethers, and salts is possible within the specific
     chemical designation, that has a degree of danger or probable
 6
 7
     danger associated with a depressant effect on the central
 8
     nervous system:
 9
          (1)
                Alprazolam;
10
          (2) Barbital;
11
          (3) Brexanolone;
12
         [\frac{(3)}{}] (4) Bromazepam;
13
         [(4)] (5) Butorphanol;
14
         [(5)] (6) Camazepam;
15
         [\frac{(6)}{(7)}] (7) Carisoprodol;
16
         [\frac{(7)}{(8)}] (8) Chloral betaine;
17
         [<del>(8)</del>] (9) Chloral hydrate;
18
         [<del>(9)</del>] (10) Chlordiazepoxide;
19
        [\frac{(10)}{(11)}] (11) Clobazam;
20
        [<del>(11)</del>] (12) Clonazepam;
21
        [<del>(12)</del>] (13) Clorazepate;
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S.B. NO. S.D. 1

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1
          [\frac{(13)}{(14)}] (14) Clotiazepam;
 2
          \left[\frac{(14)}{(15)}\right] (15) Cloxazolam;
 3
          [<del>(15)</del>] (16) Delorazepam;
 4
          [(16) Dichloralphenazone (Midrin);]
 5
           (17) Diazepam;
           (18) Dichloralphenazone (Midrin);
 6
 7
          [<del>(18)</del>] (19) Estazolam;
 8
          [\frac{(19)}{(20)}] (20) Ethchlorvynol;
 9
          \left[\frac{(20)}{(21)}\right] (21) Ethinamate;
10
          [\frac{(21)}{2}] (22) Ethyl loflazepate;
11
          [\frac{(22)}{(23)}] (23) Fludiazepam;
12
          [\frac{(23)}{(24)}] (24) Flunitrazepam;
13
          [<del>(24)</del>] (25) Flurazepam;
          [\frac{(25)}{(25)}] (26) Fospropofol (Lusedra);
14
15
          \left[\frac{(26)}{(27)}\right] (27) Halazepam;
16
          \left[\frac{(27)}{}\right] (28) Haloxazolam;
17
          \left[\frac{(28)}{(29)}\right] (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
          \left[\frac{(29)}{(31)}\right] (31) Loprazolam;
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```
1
          [<del>(30)</del>] (32) Lorazepam;
 2
          [\frac{(31)}{(33)}]
                            Lormetazepam;
 3
          [\frac{(32)}{}] (34) Mebutamate;
 4
          [\frac{(33)}{(35)}]
                            Medazepam;
 5
          [<del>(34)</del>] (36) Meprobamate;
 6
          \left[\frac{(35)}{(37)}\right] (37) Methohexital;
 7
          [<del>(36)</del>] (38)
                            Methylphenobarbital (mephorbarbital);
 8
                           Midazolam;
          [-(37)-] (39)
 9
          [\frac{(38)}{(40)}]
                           Nimetazepam;
10
          [\frac{(39)}{(41)}]
                           Nitrazepam;
11
          [\frac{(40)}{(40)}] (42) Nordiazepam;
12
          [<del>(41)</del>] (43) Oxazepam;
13
          [\frac{(42)}{(42)}] (44) Oxazolam;
14
          [<del>(43)</del>] (45) Paraldehyde;
15
          [<del>(44)</del>] (46) Petrichloral;
16
          [-(45)] (47) Phenobarbital;
17
          [<del>(46)</del>] (48) Pinazepam;
18
          \left[\frac{47}{1}\right] (49) Prazepam;
19
          [<del>(48)</del>] (50) Quazepam;
20
           (51) Remimazolam;
21
          [\frac{(49)}{(52)}]
                            Suvorexant;
```

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

```
1
         [\frac{(6)}{(7)}] (7) Mefenorex;
 2
         [\frac{7}{1}] (8) Modafinil;
 3
         [<del>(8)</del> Phentermine;
 4
          (9)
               Pemoline (including organometallic complexes and
 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}{12}] (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
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
```

S.B. NO. S.D. 1

1 central nervous system, including its salts, isomers, and salts 2 of isomers: 3 (1) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1yl]butanamide) (Other names: BRV; UCB-34714; Briviact) 4 5 and its salts; [(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 $\left[\frac{(2)}{(2)}\right]$ (4) Pregabalin $\left[\frac{(S)}{3} - \frac{(aminomethyl)}{5} - \frac{(aminomethyl)}{$ 10 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 upon its approval.

S.B. NO. 3141 S.D. 1 H.D. 2 C.D. 1

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. (CD1)

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