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: 1. By amending subsection (b) to read: 3 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
              Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-
        (12)
11
              phenethyl)-3-methyl-4-piperidinyl]-N-
12
              phenylpropanamide);
13
        (13)
              Betameprodine;
14
        (14)
              Betamethadol:
15
        (15)
              Betaprodine;
16
        (16)
              Clonitazene;
17
        (17)
              Dextromoramide;
18
        (18)
               Diampromide;
19
        (19)
               Diethylthiambutene;
20
        (20)
               Difenoxin;
21
        (21)
               Dimenoxadol;
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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
        (34)
               3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-
14
               piperidyl]-N-phenylpropanamide);
15
        (35)
               3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-
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;
6
        (45)
              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
        (55)
              Trimeperidine;
18
        (56)
              N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
19
               (benzylfentanyl), its optical isomers, salts, and
20
               salts of isomers;
```

```
1
             N-[1-(2-thienyl)methyl-4-piperidyl]-N-
        (57)
2
              phenylpropanamide (thenylfentanyl), its optical
3
              isomers, salts, and salts of isomers;
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
              isomers, esters, ethers, salts, and salts of isomers,
12
13
              esters, and ethers (Other names: Butyryl fentanyl);
14
              N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-
        (61)
15
              yl]-N-phenylpropionamide, its isomers, esters, ethers,
16
              salts and salts of isomers, esters, and ethers (Other
17
              names: beta-hydroxythiofentanyl);
18
        (62)
              N-(1-phenethylpiperidin-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);
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```
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-47700);
5
        (64)
              4-fluoroisobutyryl fentanyl or para-fluoroisobutyryl
6
              fentanyl [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-
7
              4-yl)isobutyramide];
8
        (65)
              Acryl fentanyl or acryloylfentanyl [N-(1-
9
              phenethylpiperidin-4-yl)-N-phenylacrylamide];
10
        (66)
              Ocfentanil [N-(2-fluorophenyl)-2-methoxy-N-(1-
11
              phenethylpiperidin-4-yl)acetamide];
12
        (67)
              Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
13
              phenylcyclopropanecarboxamide;
14
        (68)
              Methoxyacetyl fentanyl (2-methoxy-N-(1-
15
              phenethylpiperidin-4-yl)-N-phenylacetamide);
16
              Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-
        (69)
17
              phenethylpiperidin-4-yl)propionamide) (Other name: 2-
18
              fluorofentanyl); [and]
19
        (70)
              Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-
20
              phenethylpiperidin-4-yl)butyramide) [-];
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1	<u>(71)</u>	2'-fluoro ortho-fluorofentanyl (N-(1-(2-
2		fluorophenethyl)piperidin-4-yl)-N-(2-
3		fluorophenyl)propionamide) (Other name: 2'-fluoro 2-
4		<pre>fluorofentanyl);</pre>
5	(72)	4'-methyl acetyl fentanyl (N-(1-(4-
6		<pre>methylphenethyl)piperidin-4-yl)-N-phenylacetamide);</pre>
7	<u>(73)</u>	Beta'-phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-
8		N,3-diphenylpropanamide) (Other names: beta'-Phenyl
9		<pre>fentanyl; 3-phenylpropanoyl fentanyl);</pre>
10	(74)	Beta-methyl fentanyl (N-phenyl-N-(1-(2-
11		<pre>phenylpropyl)piperidin-4-yl)propionamide);</pre>
12	<u>(75)</u>	Ortho-fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-
13		<pre>phenethylpiperidin-4-yl)butyramide) (Other name: 2-</pre>
14		<pre>fluorobutyryl fentanyl);</pre>
15	(76)	Ortho-methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-
16		<pre>phenethylpiperidin-4-yl)acetamide) (Other name: 2-</pre>
17		<pre>methyl acetylfentanyl);</pre>
18	(77)	Ortho-methyl methoxyacetyl fentanyl (2-methoxy-N-(2-
19		methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide)
20		(Other name: 2-methyl methoxyacetyl fentanyl);

1	(78)	Para-methylfentanyl (N-(4-methylphenyl)-N-(1-
2		phenethylpiperidin-4-yl)propionamide) (Other name: 4-
3		<pre>methylfentanyl);</pre>
4	<u>(79)</u>	Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
5		<pre>phenylbenzamide) (Other name: benzoyl fentanyl);</pre>
6	(80)	Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
7		phenylthiophene-2-carboxamide) (Other names: 2-
8		thiofuranyl fentanyl; thiophene fentanyl);
9	(81)	Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-
10		<pre>yl) (phenyl) carbamate);</pre>
11	(82)	Ortho-fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-
12		<pre>phenethylpiperidin-4-yl)acrylamide);</pre>
13	(83)	Ortho-fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-
14		(1-phenethylpiperidin-4-yl)isobutyramide); and
15	(84)	Para-fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-
16		phenethylpiperidin-4-yl)furan-2-carboxamide)."
17	2.	By amending subsections (f) and (g) to read:
18	"(f)	Stimulants. Unless specifically excepted or unless
19	listed in	another schedule, any material, compound, mixture, or
20	preparati	on which contains any quantity of the following

1	substances	s having a stimulant effect on the central nervous
2	system, i	ncluding its salts, isomers, and salts of isomers:
3	(1)	Aminorex;
4	(2)	Cathinone;
5	(3)	Fenethylline;
6	(4)	Methcathinone;
7	(5)	N-ethylamphetamine;
8	(6)	4-methylaminorex;
9	(7)	N,N-dimethylamphetamine; [and]
10	(8)	Substituted cathinones, any compound, except bupropion
11		or compounds listed under a different schedule,
12		structurally derived from 2-aminopropan-1-one by
13		substitution at the 1-position with either phenyl,
14		naphthyl, or thiophene ring systems, whether or not
15		the compound is further modified in any of the
16		following ways:
17		(A) By substitution in the ring system to any extent
18		with alkyl, alkylenedioxy, alkoxy, haloalkyl,
19		hydroxyl, or halide substituents, whether or not
20		further substituted in the ring system by one or
21		more other univalent substituents;

1	(B)	By substitution at the 3-position with an acyclic
2		alkyl substituent; or
3	(C)	By substitution at the 2-amino nitrogen atom with
4		alkyl, dialkyl, benzyl, or methoxybenzyl groups,
5		or by inclusion of the 2-amino nitrogen atom in a
6		cyclic structure.
7	Some	other trade names: Mephedrone (2-methylamino-1-p-
8	toly	lpropan-1-one), also known as 4-
9	meth	ylmethcathinone (4-MMC), methylephedrone or MMCAT;
10	Meth	ylenedioxypyrovalerone (MDPV, MDPK); methylone or
11	3,4-1	methylenedioxymethcathinone; and 1-
12	(ben	zo[d][1,3]dioxol-5-yl)-2-(ethylamino)propan-1-one,
13	monol	nydrochloride, also known as Ethylone, bk-MDEA
14	hydr	ochloride, MDEC; 3,4-Methylenedioxy-N-
15	ethy	lcathinone; bk-Methylenedioxyethylamphetamine, 4-
16	meth	yl-N-ethylcathinone (4-MEC); 4-methyl-alpha-
17	pyrr	olidinopropiophenone (4-MePPP); alpha-
18	pyrr	olidinopentiophenone ([alpha]-PVP); 1-(1,3-
19	benz	odioxol-5-yl)-2-(methylamino)butan-1-one
20	(but	ylone, bk-MBDB e); 2-(methylamino)-1-phenylpentan-
21	1-on	e (pentedrone); 1-(1,3-benzodioxol-5-yl)-2-

1		(methylamino)pentan-1-one (pentylone, bk-MBDP); 4-
2		fluoro-N-methylcathinone (4-FMC, flephedrone); 3-
3		fluoro-N-methylcathinone (3-FMC); 1-(naphthalen-2-yl)-
4		2-(pyrrolidin-1-yl)pentan-1-one (naphyrone); alpha-
5		pyrrolidinobutiophenone ([alpha]-PBP) and their
6		optical, positional, and geometric isomers, salts and
7		salts of isomers, whenever the existence of such
8		salts, isomers, and salts of isomers is possible $[\div]$;
9	(9)	4,4'-dimethylaminorex (common name: 4,4'-DMAR)
10		including its salts, isomers, and salts of isomers;
11		and
12	(10)	1-(4-methoxyphenyl)-N-methylpropan-2-amine (para-
13		methoxymethamphetamine, PMMA), including its salts,
14		isomers, and salts of isomers whenever the existence
15		of such salts, isomers, and salts of isomers is
16		possible within the specific chemical designation.
17	(g)	Any of the following cannabinoids, their salts,
18	isomers,	and salts of isomers, unless specifically excepted,
19	whenever	the existence of these salts, isomers, and salts of
20	isomers i	s possible within the specific chemical designation:

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1	(1)	retranydrocannabinois; meaning tetranydrocannabinois
2		naturally contained in a plant of the genus Cannabis
3		(cannabis plant), as well as synthetic equivalents of
4		the substances contained in the plant, or in the
5		resinous extractives of Cannabis, sp. or synthetic
6		substances, derivatives, and their isomers with
7		similar chemical structure and pharmacological
8		activity to those substances contained in the plant,
9		such as the following: Delta 1 cis or trans
10		tetrahydrocannabinol, and their optical isomers; Delta
11		6 cis or trans tetrahydrocannabinol, and their optical
12		isomers; and Delta 3,4 cis or trans-
13		tetrahydrocannabinol, and its optical isomers (since
14		nomenclature of these substances is not
15		internationally standardized, compounds of these
16		structures, regardless of numerical designation of
17		atomic positions, are covered);
18	(2)	Naphthoylindoles; meaning any compound containing a 3-
19		(1-naphthoyl)indole structure with substitution at the
20		nitrogen atom of the indole ring by a alkyl,
21		haloalkyl, alkenyl, cycloalkylmethyl,cycloalkylethyl,

1		1-(N-methy1-2-piperidiny1) methy1 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 naphthyl ring to any
5		extent;
6	(3)	Naphthylmethylindoles; meaning any compound containing
7		a 1H-indol-3-yl-(1-naphthyl) methane structure with
8		substitution at the nitrogen atom of the indole ring
9		by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
10		cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or
11		2-(4-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	(4)	Naphthoylpyrroles; meaning any compound containing a
16		3-(1-naphthoyl)pyrrole structure with substitution at
17		the nitrogen atom of the pyrrole 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

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

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

```
1
        (15)
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-amino-3-methyl-1-oxobutan-2-yl)
2
               fluorobenzyl)-1H-indazole-3-carboxamide, its optical,
3
              positional, and geometric isomers, salts, and salts of
4
               isomers (Other names: AB-FUBINACA);
5
        (16)
              N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
6
               indazole-3-carboxamide, its optical, positional, and
7
               geometric isomers, salts, and salts of isomers (Other
8
               names: ADB-PINACA);
9
        (17)
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-
10
               (cyclohexylmethyl) -1H-indazole-3-carboxamide, its
11
               optical, positional, and geometric isomers, salts, and
12
               salts of isomers (Other names: AB-CHMINACA);
13
        (18)
              N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-
14
               indazole-3-carboxamide, and geometric isomers, salts,
15
               and salts of isomers (Other names: AB-PINACA);
16
        (19)
              [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-
17
               yl) methanone, and geometric isomers, salts, and salts
18
               of isomers (Other names: THJ-2201);
19
        (20)
              Methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-
20
               valinate, and geometric isomers, salts, and salts of
21
               isomers (Other names: FUB-AMB, Methyl 2-(1-(4-
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1
              fluorobenzyl)-1H-indazole-3-carboxamido)-3-
2
              methylbutanoate, MMB-FUBINACA, AMB-FUBINACA);
3
        (21)
             (S)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-
              carboxamido)-3-methylbutanoate, and geometric isomers,
4
5
              salts, and salts of isomers (Other names: 5-fluoro-
6
              AMB, 5-fluoro-AMP);
7
        (22)
             N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-
8
              indazole-3-carboxamide, and geometric isomers, salts,
9
              and salts of isomers (Other names: AKB48 N-(5-
10
              fluoropentyl) analog, 5F-AKB48, APINACA 5-fluoropentyl
11
              analog, 5F-APINACA);
12
        (23)
              N-adamantyl-1-fluoropentylindole-3-Carboxamide, and
13
              geometric isomers, salts, and salts of isomers (Other
14
              names: STS-135, 5F-APICA; 5-fluoro-APICA);
15
        (24)
             Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
16
              carboxylate, and geometric isomers, salts, and salts
17
              of isomers (Other names: NM2201);
18
        (25)
             N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-
19
              (cyclohexylmethyl) -1H-indazole-3-carboxamide, and
20
              geometric isomers, salts, and salts of isomers (Other
21
              names: MAB-CHMINACA and ADB-CHMINACA);
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1	(26)	Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-
2		carboxamido]-3,3-dimethylbutanoate (Other names: 5F-
3		ADB, 5-flouro-ADB, and 5F-MDMB-PINACA), its optical,
4		positional, and geometric isomers, salts, and salts of
5		isomers; [and]
6	(27)	1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)indazole-3-
7		carboxamide (CUMYL-4CN-BINACA), its optical,
8		positional, and geometric isomers, salts, and salts of
9		isomers; also known as SGT-78, 4-CN-CUMYL-BINACA;
10		CUMYL-CB-PINACA; CUMYL-CYBINACA; 4-cyano CUMYL-
11		BUTINACA[-];
12	(28)	Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
13		carboxylate (Other names: NM2201 or CBL2201);
14	(29)	N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-
15		fluoropentyl)-1H-indazole-3-carboxamide (Other names:
16		5F-AB-PINACA);
17	(30)	1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-
18		3-carboxamide (Other names: 4-CN-CUMYL-BUTINACA, 4-
19		cyano-CUMYL-BUTINACA; 4-CN-CUMYL BINACA, CUMYL-4CN-
20		BINACA, or SGT-78);

```
1
        (31) Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-
2
              carboxamido) - 3-methylbutanoate (Other names: MMB-
3
              CHMICA or AMB-CHMICA);
4
        (32) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-
5
              pyrrolo[2,3-b]pyridine-3-carboxamide (Other names: 5F-
6
              CUMYL-P7AICA); and
7
        (33) Methyl 3, 3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-
8
              3-carboxamido)butanoate (MDMB-4en-PINACA)."
         SECTION 2. Section 329-16, Hawaii Revised Statutes, is
9
10
    amended by amending subsection (c) to read as follows:
         "(c) Any of the following opiates, including their
11
12
    isomers, esters, ethers, salts, and salts of isomers, whenever
13
    the existence of these isomers, esters, ethers, and salts is
14
    possible within the specific chemical designation:
15
         (1)
              Alfentanil;
16
         (2)
             Alphaprodine;
17
         (3)
             Anileridine;
18
         (4)
             Bezitramide;
19
         (5)
             Bulk Dextropropoxyphene (nondosage form);
20
         (6)
             Carfentanil;
21
         (7)
             Dihydrocodeine;
```

```
1
         (8)
              Diphenoxylate;
2
         (9)
              Fentanyl;
3
              Isomethadone;
        (10)
4
        (11)
              Levo-alphacetylmethadol (LAAM);
5
        (12)
              Levomethorphan;
6
              Levorphanol;
        (13)
7
        (14)
              Metazocine;
8
        (15)
              Methadone;
9
              Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-
        (16)
10
              diphenyl butane;
11
              Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-
        (17)
12
              diphenyl-propane-carboxylic acid;
13
              Pethidine (Meperidine);
        (18)
14
        (19)
              Pethidine-Intermediate-A, 4-cyano-1-methyl-4-
15
              phenylpiperidine;
16
             Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-
        (20)
17
              carboxylate;
18
        (21)
              Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-
19
              4-carboxylic acid;
20
        (22)
              Phenazocine;
21
        (23)
             Piminodine;
```



```
1
        (24)
             Racemethorphan;
2
        (25)
             Racemorphan;
3
        (26)
             Remifentanil;
4
        (27)
             Sufentanil;
5
        (28)
             Tapentadol; [and]
6
        (29)
             Thiafentanil [-]; and
7
        (30) Oliceridine, including the free base form, and its
8
              salts, to include the fumarate salt, by definition."
9
         SECTION 3. Section 329-20, Hawaii Revised Statutes, is
10
    amended as follows:
11
         1. By amending subsection (b) to read:
12
         "(b) Depressants. Any material, compound, mixture, or
13
    preparation which contains any quantity of the following
14
    substances, including its salts, isomers, esters, ethers, and
15
    salts of isomers, whenever the existence of these isomers,
    esters, ethers, and salts is possible within the specific
16
17
    chemical designation, that has a degree of danger or probable
18
    danger associated with a depressant effect on the central
19
    nervous system:
20
             Alprazolam;
         (1)
```



(2) Barbital;

21

```
1
          (3)
               Bromazepam;
2
          (4)
               Butorphanol;
3
          (5)
               Camazepam;
4
          (6)
               Carisoprodol;
5
               Chloral betaine;
          (7)
6
          (8)
               Chloral hydrate;
7
          (9)
               Chlordiazepoxide;
8
         (10)
               Clobazam;
9
         (11)
               Clonazepam;
10
               Clorazepate;
         (12)
11
         (13)
               Clotiazepam;
12
         (14)
               Cloxazolam;
13
         (15)
               Delorazepam;
14
         (16)
               Dichloralphenazone (Midrin);
15
         (17)
               Diazepam;
16
         (18)
               Estazolam;
17
         (19)
               Ethchlorvynol;
18
         (20)
               Ethinamate;
19
               Ethyl loflazepate;
         (21)
20
         (22)
               Fludiazepam;
21
         (23)
               Flunitrazepam;
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```
Flurazepam;
1
         (24)
2
         (25)
               Fospropofol (Lusedra);
3
         (26)
               Halazepam;
4
         (27)
               Haloxazolam;
5
         (28)
               Ketazolam;
6
         (29)
               Loprazolam;
7
         (30)
               Lorazepam;
8
         (31)
               Lormetazepam;
9
         (32)
               Mebutamate;
10
               Medazepam;
         (33)
11
         (34)
               Meprobamate;
12
         (35)
               Methohexital;
13
         (36)
               Methylphenobarbital (mephorbarbital);
14
         (37)
               Midazolam;
15
         (38)
               Nimetazepam;
16
         (39)
               Nitrazepam;
17
         (40)
               Nordiazepam;
18
         (41)
               Oxazepam;
19
         (42)
              Oxazolam;
20
         (43)
               Paraldehyde;
21
         (44)
               Petrichloral;
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1
        (45)
             Phenobarbital;
2
        (46)
              Pinazepam;
3
        (47)
              Prazepam;
4
        (48)
              Quazepam;
5
        (49)
              Suvorexant;
6
        (50)
              Temazepam;
7
        (51)
              Tetrazepam;
8
        (52)
              Triazolam;
9
        (53)
              Zaleplon;
10
        (54)
              Zolpidem;
11
              Zopiclone (Lunesta); [and]
        (55)
12
        (56)
              Brexanolone [-];
13
              Remimazolam, including its salts, isomers, and salts
        (57)
14
              of isomers whenever the existence of such salts,
15
               isomers, and salts of isomers is possible; and
16
        (58) Lemborexant ((1R, 2S)-2-[(2, 4-dimethylpyrimidin-5-
17
              yl)oxymethyl]-2-(3-fluorophenyl)-N-(5-fluoropyridin-2-
18
               yl)cyclopropane-1-carboxamide), including its salts,
19
               isomers, and salts of isomers whenever the existence
20
               of such salts, isomers, and salts of isomers is
21
               possible."
```

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1
         2. By amending subsection (d) to read:
2
         "(d) Stimulants. Unless listed in another schedule, any
    material, compound, mixture, or preparation which contains any
3
4
    quantity of the following substances having a stimulant effect
    on the central nervous system, including its salts, isomers, and
5
6
    salts of such isomers whenever the existence of such salts,
7
    isomers, and salts of isomers is possible within the specific
8
    chemical designation:
9
         (1) Cathine ((+)-norpseudoephedrine);
10
             Diethylpropion;
         (2)
11
         (3)
             Fencamfamin;
12
             Fenproporex;
         (4)
13
         (5) Mazindol;
14
         (6)
             Mefenorex;
15
         (7) Modafinil;
16
         (8) Phentermine;
17
         (9)
             Pemoline (including organometallic complexes and
18
              chelates thereof);
19
        (10) Pipradrol;
20
        (11) Sibutramine;
21
        (12) SPA (1-dimethylamino-1,2-diphenylethane, lefetamine);
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1

```
(13) Lorcaserin; [and]
2
        (14) Solriamfetol [-]; and
3
        (15) Serdexmethylphenidate, including its salts, isomers,
4
              and salts of isomers."
         SECTION 4. Section 329-22, Hawaii Revised Statutes, is
5
6
    amended by amending subsection (d) to read as follows:
7
         "(d)
               Depressants. Unless specifically exempted or
8
    excluded or unless listed in another schedule, any material,
9
    compound, mixture, or preparation that contains any quantity of
    the following substances having a depressant effect on the
10
11
    central nervous system, including its salts, isomers, and salts
    of isomers:
12
13
              Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-
         (1)
14
              propionamide], (Vimpat);
15
         (2)
             Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic
16
              acid]; [and]
17
         (3)
             Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-
18
              yl]butanamide) (Other names: BRV; UCB-34714; Briviact)
19
              and its salts[-]; and
20
         (4) Lasmiditan (2,4,6-trifluoro-N-(6-(1-methylpiperidine-
21
              4-carbonyl)pyridine-2-yl-benzamide)."
```

- 1 SECTION 5. Statutory material to be repealed is bracketed
- 2 and stricken. New statutory material is underscored.
- 3 SECTION 6. This Act shall take effect on January 1, 2050.

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. (HD1)

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