

**ACT 233**

H.B. NO. 1032

A Bill for an Act Relating to the Uniform Controlled Substances Act.

*Be It Enacted by the Legislature of the State of Hawaii:*

SECTION 1. Section 329-14, Hawaii Revised Statutes, is amended as follows:

1. By amending subsection (b) to read:

“(b) Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

- (1) Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);
- (2) Acetylmethadol;
- (3) Allylprodine;
- (4) Alphacetylmethadol (except levo-alphacetylmethadol, levomethadyl acetate, or LAAM);
- (5) Alphameprodine;
- (6) Alphamethadol;
- (7) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl] propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);
- (8) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
- (9) Benzethidine;
- (10) Betacetylmethadol;
- (11) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide);
- (12) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide);
- (13) Betameprodine;
- (14) Betamethadol;
- (15) Betaprodine;
- (16) Clonitazene;
- (17) Dextromoramide;
- (18) Diampromide;
- (19) Diethylthiambutene;
- (20) Difenoxyin;
- (21) Dimenoxadol;
- (22) Dimepheptanol;
- (23) Dimethylthiambutene;
- (24) Dioxaphetyl butyrate;
- (25) Dipipanone;
- (26) Ethylmethylthiambutene;
- (27) Etonitazene;
- (28) Etoxidine;
- (29) Furethidine;
- (30) Hydroxypethidine;
- (31) Ketobemidone;
- (32) Levomoramide;
- (33) Levophenacilmorphan;
- (34) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);
- (35) 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
- (36) Morpheridine;
- (37) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- (38) Noracymethadol;
- (39) Norlevorphanol;
- (40) Normethadone;

- (41) Norpipanone;
- (42) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide;
- (43) PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- (44) Phenadoxone;
- (45) Phenampromide;
- (46) Phenomorphan;
- (47) Phenoperidine;
- (48) Piritramide;
- (49) Proheptazine;
- (50) Properidine;
- (51) Propiram;
- (52) Racemoramide;
- (53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);
- (54) Tilidine;
- (55) Trimeperidine;
- (56) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts, and salts of isomers;
- (57) N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical isomers, salts, and salts of isomers;
- (58) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide, (acetyl fentanyl), its optical, positional, and geometric isomers, salts, and salts of isomers;
- (59) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)cyclohexylmethyl]benzamide), its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers;
- (60) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: Butyryl fentanyl);
- (61) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: beta-hydroxythiofentanyl);
- (62) N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: Furanyl fentanyl);
- (63) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: U-47700);
- (64) 4-fluoroisobutyryl fentanyl or para-fluoroisobutyryl [~~fentanyl~~ fentanyl] [N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide];
- (65) Acryl fentanyl or acryloylfentanyl [N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]; [~~and~~]
- (66) Ocfentanil [~~[N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide]-] [N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide];~~
- (67) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide;
- (68) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

- (69) Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide) (Other name: 2-fluorofentanyl); and
- (70) Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide).”
2. By amending subsection (e) to read:
- “(e) Depressants. Unless specifically excepted, the schedule shall include any material, compound, mixture, or preparation which contains any quantity of the substance:
- (1) Mecloqualone; [øƒ]
  - (2) Methaqualone[-];
  - (3) Etizolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible); or
  - (4) Flualprazolam (including its optical, positional, and geometric isomers, salts, and salts of isomers, where possible).”
3. By amending subsection (g) to read:
- “(g) Any of the following cannabinoids, their salts, isomers, and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:
- (1) Tetrahydrocannabinols; meaning tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. or synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following: Delta 1 cis or trans tetrahydrocannabinol, and their optical isomers; Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers; and Delta 3,4 cis or trans-tetrahydrocannabinol, and its optical isomers (since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions, are covered);
  - (2) Naphthoylindoles; meaning any compound containing a 3-(1-naphthoyl)indole structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent;
  - (3) Naphthylmethylindoles; meaning any compound containing a 1H-indol-3-yl-(1-naphthyl) methane structure with substitution at the nitrogen atom of the indole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent;
  - (4) Naphthoylpyrroles; meaning any compound containing a 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom of the pyrrole ring by a alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl) ethyl group whether or not further substituted in the pyrrole ring to any extent, whether or not substituted in the naphthyl ring to any extent;

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

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

SECTION 2. Section 329-16, Hawaii Revised Statutes, is amended by amending subsection (f) to read as follows:

“(f) Immediate precursor. Unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

- (1) Immediate precursor to amphetamine and methamphetamine:

- (A) Phenylacetone, phenyl-2-propanone(P2P), benzyl methyl ketone, methyl benzyl ketone;
- (2) Immediate precursors to phencyclidine (PCP):
  - (A) 1-phenylcyclohexylamine; ~~and~~ or
  - (B) 1-piperidinocyclohexanecarbonitrile(PCC); or
- (3) Immediate precursor to Fentanyl:
  - (A) 4-anilino-N-phenethyl-4-piperidine (ANPP)[-]; or
  - (B) N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl).”

SECTION 3. Section 329-20, Hawaii Revised Statutes, is amended as follows:

1. By amending subsection (b) to read:

“(b) Depressants. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, esters, ethers, and salts of isomers, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation, that has a degree of danger or probable danger associated with a depressant effect on the central nervous system:

- (1) Alprazolam;
- (2) Barbitol;
- (3) Bromazepam;
- (4) Butorphanol;
- (5) Camazepam;
- (6) Carisoprodol;
- (7) Chloral betaine;
- (8) Chloral hydrate;
- (9) Chlordiazepoxide;
- (10) Clobazam;
- (11) Clonazepam;
- (12) Clorazepate;
- (13) Clotiazepam;
- (14) Cloxazolam;
- (15) Delorazepam;
- (16) Dichloralphenazone (Midrin);
- (17) Diazepam;
- (18) Estazolam;
- (19) Ethchlorvynol;
- (20) Ethinamate;
- (21) Ethyl loflazepate;
- (22) Fludiazepam;
- (23) Flunitrazepam;
- (24) Flurazepam;
- (25) Fospropofol (Lusedra);
- (26) Halazepam;
- (27) Haloxazolam;
- (28) Ketazolam;
- (29) Loprazolam;
- (30) Lorazepam;
- (31) Lormetazepam;
- (32) Mebutamate;
- (33) Medazepam;
- (34) Meprobamate;
- (35) Methohexital;
- (36) Methylphenobarbital (mephorbarbital);

- (37) Midazolam;
- (38) Nimetazepam;
- (39) Nitrazepam;
- (40) Nordiazepam;
- (41) Oxazepam;
- (42) Oxazolam;
- (43) Paraldehyde;
- (44) Petrichloral;
- (45) Phenobarbital;
- (46) Pinazepam;
- (47) Prazepam;
- (48) Quazepam;
- (49) Suvorexant;
- (50) Temazepam;
- (51) Tetrazepam;
- (52) Triazolam;
- (53) Zaleplon;
- (54) Zolpidem; [~~and~~]
- (55) Zopiclone (Lunesta)[-]; and
- (56) Brexanolone.”

2. By amending subsection (d) to read:

“(d) Stimulants. Unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Cathine ((+)-norpseudoephedrine);
- (2) Diethylpropion;
- (3) Fencamfamin;
- (4) Fenproporex;
- (5) Mazindol;
- (6) Mefenorex;
- (7) Modafinil;
- (8) Phentermine;
- (9) Pemoline (including organometallic complexes and chelates thereof);
- (10) Pipradrol;
- (11) Sibutramine;
- (12) SPA (1-dimethylamino-1,2-diphenylethane, lefetamine); [~~and~~]
- (13) Lorcaserin[-]; and
- (14) Solriamfetol.”

SECTION 4. Section 329-22, Hawaii Revised Statutes, is amended to read as follows:

“**§329-22 Schedule V.** (a) The controlled substances listed in this section are included in schedule V.

(b) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs, which also contains one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation, valuable medicinal qualities other than those possessed by the narcotic drug alone:

- (1) Not more than 200 milligrams of codeine, or any of its salts, per 100 milliliters or per 100 grams;



- (2) Not more than 100 milligrams of dihydrocodeine, or any of its salts, per 100 milliliters or per 100 grams;
- (3) Not more than 100 milligrams of ethylmorphine, or any of its salts, per 100 milliliters or per 100 grams;
- (4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit;
- (5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams; and
- (6) Not more than 0.5 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.

(c) Stimulants. Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation that contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of isomers.

(d) Depressants. Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation that contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers:

- (1) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxypropionamide], (Vimpat);
- (2) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid]; and
- (3) Brivaracetam ((2S)-2'[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (Other names: BRV; UCB-234714; Briviact) and its salts.

~~[(e) Approved cannabidiol drugs. A drug product in finished dosage formulation that has been approved by the United States Food and Drug Administration that contains cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-1,3-benzenediol) derived from cannabis and no more than 0.1 percent<sup>3</sup> (w/w) residual tetrahydrocannabinols.]”~~

SECTION 5. Statutory material to be repealed is bracketed and stricken. New statutory material is underscored.

SECTION 6. This Act shall take effect upon its approval.

(Approved July 6, 2021.)

**Notes**

- 1. Prior to amendment a hyphen appeared here.
- 2. Prior to amendment a dash appeared here.
- 3. Prior to amendment “per cent” appeared here.