
SUBSTITUTE SENATE BILL 5101

State of Washington

62nd Legislature

2011 Regular Session

By Senate Judiciary (originally sponsored by Senators Carrell, Schoesler, Delvin, and Stevens)

READ FIRST TIME 02/10/11.

1 AN ACT Relating to placing certain synthetic cannabimimetics into
2 schedule I of the uniform controlled substances act; amending RCW
3 69.50.204; creating a new section; and declaring an emergency.

4 BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF WASHINGTON:

5 NEW SECTION. **Sec. 1.** The legislature finds that synthetic
6 cannabimimetics have been developed for research purposes to
7 investigate the cannabimimetic system. No legitimate nonresearch uses
8 have been identified for synthetic cannabimimetics nor have they been
9 approved for human consumption by the United States food and drug
10 administration.

11 The legislature further finds that the popularity of synthetic
12 cannabimimetics has greatly increased in the United States and they are
13 being abused for their psychoactive properties. Products containing
14 synthetic cannabimimetics are marketed as legal alternatives to
15 marijuana and are being sold over the internet and in tobacco and smoke
16 shops, drug paraphernalia shops, and convenience stores. Due to their
17 method of manufacture and high pharmacological potency, synthetic
18 cannabimimetics are potentially extremely harmful; for some users,
19 smoking synthetic cannabimimetics for the purpose of achieving

1 intoxication and experiencing the psychoactive effects has led to
2 emergency room visits and calls to poison control centers.
3 Consequently, placement of the synthetic cannabimimetics listed in RCW
4 69.50.204(c)(30) into schedule I of the uniform controlled substances
5 act is necessary to avoid an imminent hazard to public safety.

6 **Sec. 2.** RCW 69.50.204 and 2010 c 177 s 2 are each amended to read
7 as follows:

8 Unless specifically excepted by state or federal law or regulation
9 or more specifically included in another schedule, the following
10 controlled substances are listed in Schedule I:

11 (a) Any of the following opiates, including their isomers, esters,
12 ethers, salts, and salts of isomers, esters, and ethers whenever the
13 existence of these isomers, esters, ethers, and salts is possible
14 within the specific chemical designation:

15 (1) Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-
16 piperidinyl]-N-phenylacetamide);

17 (2) Acetylmethadol;

18 (3) Allylprodine;

19 (4) Alphacetylmethadol, except levo-alphacetylmethadol, also known
20 as levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;

21 (5) Alphameprodine;

22 (6) Alphamethadol;

23 (7) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl) ethyl-
24 4-piperidyl] propionanilide); (1-(1-methyl-2-phenylethyl)-4-(N-
25 propanilido) piperidine);

26 (8) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-
27 piperidinyl]-N-phenylpropanamide);

28 (9) Benzethidine;

29 (10) Betacetylmethadol;

30 (11) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-
31 piperidinyl]-N-phenylpropanamide);

32 (12) Beta-hydroxy-3-methylfentanyl, some trade or other names: N-
33 [1-(2-hydrox-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;

34 (13) Betameprodine;

35 (14) Betamethadol;

36 (15) Betaprodine;

37 (16) Clonitazene;

- 1 (17) Dextromoramide;
- 2 (18) Diampromide;
- 3 (19) Diethylthiambutene;
- 4 (20) Difenoxin;
- 5 (21) Dimenoxadol;
- 6 (22) Dimepheptanol;
- 7 (23) Dimethylthiambutene;
- 8 (24) Dioxaphetyl butyrate;
- 9 (25) Dipipanone;
- 10 (26) Ethylmethylthiambutene;
- 11 (27) Etonitazene;
- 12 (28) Etoxeridine;
- 13 (29) Furethidine;
- 14 (30) Hydroxypethidine;
- 15 (31) Ketobemidone;
- 16 (32) Levomoramide;
- 17 (33) Levophenacymorphan;
- 18 (34) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-
- 19 N-phenylprop anamide);
- 20 (35) 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
- 21 piperidinyl]-N-phenylpropanamide);
- 22 (36) Morpheridine;
- 23 (37) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- 24 (38) Noracymethadol;
- 25 (39) Norlevorphanol;
- 26 (40) Normethadone;
- 27 (41) Norpipanone;
- 28 (42) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-
- 29 piperidinyl] propanamide);
- 30 (43) PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 31 (44) Phenadoxone;
- 32 (45) Phenampromide;
- 33 (46) Phenomorphan;
- 34 (47) Phenoperidine;
- 35 (48) Piritramide;
- 36 (49) Proheptazine;
- 37 (50) Properidine;
- 38 (51) Propiram;

- 1 (52) Racemoramide;
2 (53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-
3 propanamide);
4 (54) Tilidine;
5 (55) Trimeperidine.

6 (b) Opium derivatives. Unless specifically excepted or unless
7 listed in another schedule, any of the following opium derivatives,
8 including their salts, isomers, and salts of isomers whenever the
9 existence of those salts, isomers, and salts of isomers is possible
10 within the specific chemical designation:

- 11 (1) Acetorphine;
12 (2) Acetyldihydrocodeine;
13 (3) Benzylmorphine;
14 (4) Codeine methylbromide;
15 (5) Codeine-N-Oxide;
16 (6) Cyprenorphine;
17 (7) Desomorphine;
18 (8) Dihydromorphine;
19 (9) Drotebanol;
20 (10) Etorphine, except hydrochloride salt;
21 (11) Heroin;
22 (12) Hydromorphenol;
23 (13) Methyldesorphine;
24 (14) Methyldihydromorphine;
25 (15) Morphine methylbromide;
26 (16) Morphine methylsulfonate;
27 (17) Morphine-N-Oxide;
28 (18) Myrophine;
29 (19) Nicocodeine;
30 (20) Nicomorphine;
31 (21) Normorphine;
32 (22) Pholcodine;
33 (23) Thebacon.

34 (c) Hallucinogenic substances. Unless specifically excepted or
35 unless listed in another schedule, any material, compound, mixture, or
36 preparation which contains any quantity of the following hallucinogenic
37 substances, including their salts, isomers, and salts of isomers
38 whenever the existence of those salts, isomers, and salts of isomers is

1 possible within the specific chemical designation. For the purposes of
2 this subsection only, the term "isomer" includes the optical, position,
3 and geometric isomers:

4 (1) Alpha-ethyltryptamine: Some trade or other names:
5 Etryptamine; monase; α -ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)
6 indole; α -ET; and AET;

7 (2) 4-bromo-2,5-dimethoxy-amphetamine: Some trade or other names:
8 4-bromo-2,5-dimethoxy- α -methylphenethylamine; 4-bromo-2,5-DMA;

9 (3) 4-bromo-2,5-dimethoxyphenethylamine: Some trade or other
10 names: 2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl
11 DOB; 2C-B, nexus;

12 (4) 2,5-dimethoxyamphetamine: Some trade or other names: 2,5-
13 dimethoxy- α -methylphenethylamine; 2,5-DMA;

14 (5) 2,5-dimethoxy-4-ethylamphetamine (DOET);

15 (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine: Other name:
16 2C-T-7;

17 (7) 4-methoxyamphetamine: Some trade or other names: 4-methoxy- α -
18 methylphenethylamine; paramethoxyamphetamine, PMA;

19 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;

20 (9) 4-methyl-2,5-dimethoxy-amphetamine: Some trade and other
21 names: 4-methyl-2,5-dimethoxy- α -methylphenethylamine; "DOM"; and
22 "STP";

23 (10) 3,4-methylenedioxy amphetamine;

24 (11) 3,4-methylenedioxymethamphetamine (MDMA);

25 (12) 3,4-methylenedioxy-N-ethylamphetamine, also known as N-ethyl-
26 alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA;

27 (13) N-hydroxy-3,4-methylenedioxyamphetamine also known as
28 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, N-hydroxy MDA;

29 (14) 3,4,5-trimethoxy amphetamine;

30 (15) Alpha-methyltryptamine: Other name: AMT;

31 (16) Bufotenine: Some trade or other names: 3-(beta-
32 Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol;
33 N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;

34 (17) Diethyltryptamine: Some trade or other names: N,N-
35 Diethyltryptamine; DET;

36 (18) Dimethyltryptamine: Some trade or other names: DMT;

37 (19) 5-methoxy-N,N-diisopropyltryptamine: Other name: 5-MeO-DIPT;

- 1 (20) Ibogaine: Some trade or other names: 7-Ethyl-6,6
2 beta,7,8,9,10,12,13,-octahydro-2-methoxy-6,9-methano-5H-pyrido (1',2'
3 1,2) azepino (5,4-b) indole; Tabernanthe iboga;
- 4 (21) Lysergic acid diethylamide;
- 5 (22) Marihuana or marijuana;
- 6 (23) Mescaline;
- 7 (24) Parahexyl-7374: Some trade or other names: 3-Hexyl-1-
8 hydroxy-7, 8, 9, 10-tetrahydro-6, 6, 9-trimethyl-6H-dibenzo[b,d]pyran;
9 synhexyl;
- 10 (25) Peyote, meaning all parts of the plant presently classified
11 botanically as Lophophora Williamsii Lemaire, whether growing or not,
12 the seeds thereof, any extract from any part of such plant, and every
13 compound, manufacture, salts, derivative, mixture, or preparation of
14 such plant, its seeds, or extracts; (interprets 21 U.S.C. Sec. 812 (c),
15 Schedule I (c)(12));
- 16 (26) N-ethyl-3-piperidyl benzilate;
- 17 (27) N-methyl-3-piperidyl benzilate;
- 18 (28) Psilocybin;
- 19 (29) Psilocyn;
- 20 (30) Any of the following synthetic cannabimimetics, their salts,
21 isomers, and salts of isomers, unless specifically excepted, whenever
22 the existence of these salts, isomers, and salts of isomers is possible
23 within the specific chemical designation:
- 24 (i) Naphthoylindoles: Any compound containing a
25 3-(1-naphthoyl)indole structure with substitution at the nitrogen atom
26 of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
27 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
28 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
29 indole ring to any extent and whether or not substituted in the
30 naphthyl ring to any extent;
- 31 (ii) Naphthylmethylindoles: Any compound containing a
32 1H-indol-3-yl-(1-naphthyl)methane structure with substitution at the
33 nitrogen atom of the indole ring by an alkyl, haloalkyl, alkenyl,
34 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
35 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
36 indole ring to any extent and whether or not substituted in the
37 naphthyl ring to any extent;

1 (iii) Naphthoylpyrroles: Any compound containing a
2 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom
3 of the pyrrole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
4 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
5 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
6 pyrrole ring to any extent and whether or not substituted in the
7 naphthyl ring to any extent;

8 (iv) Naphthylmethylindenes: Any compound containing a
9 naphthylideneindene structure with substitution at the 3-position of
10 the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
11 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
12 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
13 indene ring to any extent and whether or not substituted in the
14 naphthyl ring to any extent;

15 (v) Phenylacetylindoles: Any compound containing a
16 3-phenylacetylindole structure with substitution at the nitrogen atom
17 of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
18 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
19 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
20 indole ring to any extent and whether or not substituted in the phenyl
21 ring to any extent;

22 (vi) Cyclohexylphenols: Any compound containing a
23 2-(3-hydroxycyclohexyl)phenol structure with substitution at the
24 5-position of the phenolic ring by an alkyl, haloalkyl, alkenyl,
25 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
26 2-(4-morpholinyl)ethyl group, whether or not substituted in the
27 cyclohexyl ring to any extent;

28 (vii) Benzoylindoles: Any compound containing a 3-(benzoyl)indole
29 structure with substitution at the nitrogen atom of the indole ring by
30 an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
31 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group,
32 whether or not further substituted in the indole ring to any extent and
33 whether or not substituted in the phenyl ring to any extent;

34 (viii) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo
35 [1,2,3-de]-1,4-benzoxazin-6-yl]-1-naphthalenylmethanone: Some trade or
36 other names: WIN 55,212-2;

37 (31) Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally
38 contained in a plant of the genus Cannabis (cannabis plant), as well as

1 synthetic equivalents of the substances contained in the plant, or in
2 the resinous extractives of Cannabis, species, and/or synthetic
3 substances, derivatives, and their isomers with similar chemical
4 structure and pharmacological activity such as the following:

5 (i) 1 - cis - or trans tetrahydrocannabinol, and their optical
6 isomers, excluding tetrahydrocannabinol in sesame oil and encapsulated
7 in a soft gelatin capsule in a drug product approved by the United
8 States Food and Drug Administration;

9 (ii) 6 - cis - or trans tetrahydrocannabinol, and their optical
10 isomers;

11 (iii) 3,4 - cis - or trans tetrahydrocannabinol, and its optical
12 isomers;

13 (iv) (6aR,10aR)-9-(hydroxymethyl)-6, 6-dimethyl-3-
14 (2-methyloctan-2-yl)-6a,7,10, 10a-tetrahydrobenzo[c]chromen-1-ol: Some
15 trade or other names: HU-210;

16 (Since nomenclature of these substances is not internationally
17 standardized, compounds of these structures, regardless of numerical
18 designation of atomic positions covered.)

19 ~~((+31))~~ (32) Ethylamine analog of phencyclidine: Some trade or
20 other names: N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)
21 ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE;

22 ~~((+32))~~ (33) Pyrrolidine analog of phencyclidine: Some trade or
23 other names: 1-(1-phenylcyclohexyl)pyrrolidine; PCPy; PHP;

24 ~~((+33))~~ (34) Thiophene analog of phencyclidine: Some trade or
25 other names: 1-(1-[2-thienyl]-cyclohexyl)-piperidine; 2-thienyl analog of
26 phencyclidine; TPCP; TCP;

27 ~~((+34))~~ (35) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine: A trade or
28 other name is TCPy.

29 (d) Depressants. Unless specifically excepted or unless listed in
30 another schedule, any material, compound, mixture, or preparation which
31 contains any quantity of the following substances having a depressant
32 effect on the central nervous system, including its salts, isomers, and
33 salts of isomers whenever the existence of such salts, isomers, and
34 salts of isomers is possible within the specific chemical designation.

35 (1) Gamma-hydroxybutyric acid: Some other names include GHB;
36 gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic acid;
37 sodium oxybate; sodium oxybutyrate;

38 (2) Mecloqualone;

1 (3) Methaqualone.

2 (e) Stimulants. Unless specifically excepted or unless listed in
3 another schedule, any material, compound, mixture, or preparation which
4 contains any quantity of the following substances having a stimulant
5 effect on the central nervous system, including its salts, isomers, and
6 salts of isomers:

7 (1) Aminorex: Some other names: aminoxaphen; 2-amino-5-phenyl-2-
8 oxazoline; or 4, 5-dihydro-5-phenly-2-oxazolamine;

9 (2) N-Benzylpiperazine: Some other names: BZP, 1-benzylpiperazine;

10 (3) Cathinone, also known as 2-amino-1-phenyl-1-propanone,
11 alpha-aminopropiophenone, 2-aminopropiophenone and norephedrone;

12 (4) Fenethylamine;

13 (5) Methcathinone: Some other names: 2-(methylamino)-
14 propiophenone; alpha-(methylamino)propiophenone; 2-(methylamino)-1-
15 phenylpropan-1-one; alpha-N-methylaminopropiophenone;
16 monomethylpropion; ephedrone; N-methylcathinone; methylcathinone; AL-
17 464; AL-422; AL-463 and UR1432, its salts, optical isomers, and salts
18 of optical isomers;

19 (6) (+-)-cis-4-methylaminorex ((+)-cis-4,5-dihydro-4-methyl-5-
20 phenyl-2-oxazolamine);

21 (7) N-ethylamphetamine;

22 (8) N,N-dimethylamphetamine: Some trade or other names: N,N-
23 alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenoethylene.

24 The controlled substances in this section may be added,
25 rescheduled, or deleted as provided for in RCW 69.50.201.

26 NEW SECTION. **Sec. 3.** This act is necessary for the immediate
27 preservation of the public peace, health, or safety, or support of the
28 state government and its existing public institutions, and takes effect
29 immediately.

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