
SENATE BILL 6007

State of Washington 62nd Legislature 2012 Regular Session

By Senators Carrell, Schoesler, Benton, Regala, Delvin, and Parlette

Read first time 01/09/12. Referred to Committee on Judiciary.

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

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

6 NEW SECTION. **Sec. 1.** The legislature finds that synthetic
7 substances have been developed for research purposes. Legitimate
8 nonresearch uses have neither been identified for synthetic
9 cannabimimetics or certain substituted cathinones nor have they been
10 approved for human consumption by the United States food and drug
11 administration.

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

1 extremely harmful; for some users, smoking synthetic cannabimimetics
2 for the purpose of achieving intoxication and experiencing the
3 psychoactive effects has led to emergency room visits and calls to
4 poison control centers. Certain substituted cathinones are similarly
5 being deceptively marketed as bath salts, and have been abused by
6 persons in the state in order to achieve intoxication, resulting in
7 emergency calls and hospitalizations. Consequently, placement of the
8 synthetic cannabimimetics listed in RCW 69.50.204(c)(30) and certain
9 substituted cathinones into schedule I of the uniform controlled
10 substances act is necessary to avoid an imminent hazard to public
11 safety.

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

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

17 (a) Any of the following opiates, including their isomers, esters,
18 ethers, salts, and salts of isomers, esters, and ethers whenever the
19 existence of these isomers, esters, ethers, and salts is possible
20 within the specific chemical designation:

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

23 (2) Acetylmethadol;

24 (3) Allylprodine;

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

27 (5) Alphameprodine;

28 (6) Alphamethadol;

29 (7) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl) ethyl-
30 4-piperidyl] propionanilide); (1-(1-methyl-2-phenylethyl)-4-(N-
31 propanilido) piperidine);

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

34 (9) Benzethidine;

35 (10) Betacetylmethadol;

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

- 1 (12) Beta-hydroxy-3-methylfentanyl, some trade or other names: N-
- 2 [1-(2-hydrox-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;
- 3 (13) Betameprodine;
- 4 (14) Betamethadol;
- 5 (15) Betaprodine;
- 6 (16) Clonitazene;
- 7 (17) Dextromoramide;
- 8 (18) Diampromide;
- 9 (19) Diethylthiambutene;
- 10 (20) Difenoquin;
- 11 (21) Dimenoxadol;
- 12 (22) Dimepheptanol;
- 13 (23) Dimethylthiambutene;
- 14 (24) Dioxaphetyl butyrate;
- 15 (25) Dipipanone;
- 16 (26) Ethylmethylthiambutene;
- 17 (27) Etonitazene;
- 18 (28) Etoxadine;
- 19 (29) Furethidine;
- 20 (30) Hydroxypethidine;
- 21 (31) Ketobemidone;
- 22 (32) Levomoramide;
- 23 (33) Levophenacymorphan;
- 24 (34) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-
- 25 N-phenylprop anamide);
- 26 (35) 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-
- 27 piperidinyl]-N-phenylpropanamide);
- 28 (36) Morpheridine;
- 29 (37) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- 30 (38) Noracymethadol;
- 31 (39) Norlevorphanol;
- 32 (40) Normethadone;
- 33 (41) Norpipanone;
- 34 (42) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-
- 35 piperidinyl] propanamide);
- 36 (43) PEPAP(1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 37 (44) Phenadoxone;
- 38 (45) Phenampromide;

- 1 (46) Phenomorphan;
- 2 (47) Phenoperidine;
- 3 (48) Piritramide;
- 4 (49) Proheptazine;
- 5 (50) Properidine;
- 6 (51) Propiram;
- 7 (52) Racemoramide;
- 8 (53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-
- 9 propanamidine);
- 10 (54) Tilidine;
- 11 (55) Trimeperidine.

12 (b) Opium derivatives. Unless specifically excepted or unless
13 listed in another schedule, any of the following opium derivatives,
14 including their salts, isomers, and salts of isomers whenever the
15 existence of those salts, isomers, and salts of isomers is possible
16 within the specific chemical designation:

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

1 (23) Thebacon.

2 (c) Hallucinogenic substances. Unless specifically excepted or
3 unless listed in another schedule, any material, compound, mixture, or
4 preparation which contains any quantity of the following hallucinogenic
5 substances, including their salts, isomers, and salts of isomers
6 whenever the existence of those salts, isomers, and salts of isomers is
7 possible within the specific chemical designation. For the purposes of
8 this subsection only, the term "isomer" includes the optical, position,
9 and geometric isomers:

10 (1) Alpha-ethyltryptamine: Some trade or other names:
11 Etryptamine; monase; a-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)
12 indole; a-ET; and AET;

13 (2) 4-bromo-2,5-dimethoxy-amphetamine: Some trade or other names:
14 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA;

15 (3) 4-bromo-2,5-dimethoxyphenethylamine: Some trade or other
16 names: 2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl
17 DOB; 2C-B, nexus;

18 (4) 2,5-dimethoxyamphetamine: Some trade or other names: 2,5-
19 dimethoxy-a-methylphenethylamine; 2,5-DMA;

20 (5) 2,5-dimethoxy-4-ethylamphetamine (DOET);

21 (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine: Other name:
22 2C-T-7;

23 (7) 4-methoxyamphetamine: Some trade or other names: 4-methoxy-a-
24 methylphenethylamine; paramethoxyamphetamine, PMA;

25 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;

26 (9) 4-methyl-2,5-dimethoxy-amphetamine: Some trade and other
27 names: 4-methyl-2,5-dimethoxy-a-methylphenethylamine; "DOM"; and
28 "STP";

29 (10) 3,4-methylenedioxy amphetamine;

30 (11) 3,4-methylenedioxymethamphetamine (MDMA);

31 (12) 3,4-methylenedioxy-N-ethylamphetamine, also known as N-ethyl-
32 alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA;

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

35 (14) 3,4,5-trimethoxy amphetamine;

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

37 (16) Bufotenine: Some trade or other names: 3-(beta-

- 1 Dimethylaminoethyl)-5-hydroxindole; 3-(2-dimethylaminoethyl)-5-indolol;
2 N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;
3 (17) Diethyltryptamine: Some trade or other names: N,N-
4 Diethyltryptamine; DET;
5 (18) Dimethyltryptamine: Some trade or other names: DMT;
6 (19) 5-methoxy-N,N-diisopropyltryptamine: Other name: 5-MeO-DIPT;
7 (20) Ibogaine: Some trade or other names: 7-Ethyl-6,6
8 beta,7,8,9,10,12,13,-octahydro-2-methoxy-6,9-methano-5H-pyndo (1',2'
9 1,2) azepino (5,4-b) indole; Tabernanthe iboga;
10 (21) Lysergic acid diethylamide;
11 (22) Marihuana or marijuana;
12 (23) Mescaline;
13 (24) Parahexyl-7374: Some trade or other names: 3-Hexyl-1-
14 hydroxy-7, 8, 9, 10-tetrahydro-6, 6, 9-trimethyl-6H-dibenzo[b,d]pyran;
15 synhexyl;
16 (25) Peyote, meaning all parts of the plant presently classified
17 botanically as Lophophora Williamsii Lemaire, whether growing or not,
18 the seeds thereof, any extract from any part of such plant, and every
19 compound, manufacture, salts, derivative, mixture, or preparation of
20 such plant, its seeds, or extracts; (interprets 21 U.S.C. Sec. 812 (c),
21 Schedule I (c)(12));
22 (26) N-ethyl-3-piperidyl benzilate;
23 (27) N-methyl-3-piperidyl benzilate;
24 (28) Psilocybin;
25 (29) Psilocyn;
26 (30) Any of the following synthetic cannabimimetics, their salts,
27 isomers, and salts of isomers, unless specifically excepted, whenever
28 the existence of these salts, isomers, and salts of isomers is possible
29 within the specific chemical designation:
30 (i) Naphthoylindoles: Any compound containing a
31 3-(1-naphthoyl)indole structure with substitution at the nitrogen atom
32 of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
33 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
34 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
35 indole ring to any extent and whether or not substituted in the
36 naphthyl ring to any extent;
37 (ii) Naphthylmethylinindoles: Any compound containing a
38 1H-indol-3-yl-(1-naphthyl)methane structure with substitution at the

1 nitrogen atom of the indole ring by an alkyl, haloalkyl, alkenyl,
2 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
3 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
4 indole ring to any extent and whether or not substituted in the
5 naphthyl ring to any extent;

6 (iii) Naphthoylpyrroles: Any compound containing a
7 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom
8 of the pyrrole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
9 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
10 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
11 pyrrole ring to any extent and whether or not substituted in the
12 naphthyl ring to any extent;

13 (iv) Naphthylmethylindenes: Any compound containing a
14 naphthylideneindene structure with substitution at the 3-position of
15 the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
16 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
17 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
18 indene ring to any extent and whether or not substituted in the
19 naphthyl ring to any extent;

20 (v) Phenylacetylindoles: Any compound containing a
21 3-phenylacetylindole structure with substitution at the nitrogen atom
22 of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl,
23 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
24 2-(4-morpholinyl)ethyl group, whether or not further substituted in the
25 indole ring to any extent and whether or not substituted in the phenyl
26 ring to any extent;

27 (vi) Cyclohexylphenols: Any compound containing a
28 2-(3-hydroxycyclohexyl)phenol structure with substitution at the
29 5-position of the phenolic ring by an alkyl, haloalkyl, alkenyl,
30 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or
31 2-(4-morpholinyl)ethyl group, whether or not substituted in the
32 cyclohexyl ring to any extent;

33 (vii) Benzoylindoles: Any compound containing a 3-(benzoyl)indole
34 structure with substitution at the nitrogen atom of the indole ring by
35 an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
36 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group,
37 whether or not further substituted in the indole ring to any extent and
38 whether or not substituted in the phenyl ring to any extent;

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

4 (31) Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally
5 contained in a plant of the genus Cannabis (cannabis plant), as well as
6 synthetic equivalents of the substances contained in the plant, or in
7 the resinous extractives of Cannabis, species, and/or synthetic
8 substances, derivatives, and their isomers with similar chemical
9 structure and pharmacological activity such as the following:

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

14 (ii) 6 - cis - or trans tetrahydrocannabinol, and their optical
15 isomers;

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

18 (iv) (6aR,10aR)-9-(hydroxymethyl)-6, 6-dimethyl-3-
19 (2-methyloctan-2-yl)-6a,7,10, 10a-tetrahydrobenzo[c]chromen-1-ol: Some
20 trade or other names: HU-210;

21 (Since nomenclature of these substances is not internationally
22 standardized, compounds of these structures, regardless of numerical
23 designation of atomic positions covered.)

24 ~~((+31))~~ (32) Ethylamine analog of phencyclidine: Some trade or
25 other names: N-ethyl-1phenylcyclohexylamine, (1-phenylcyclohexyl)
26 ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE;

27 ~~((+32))~~ (33) Pyrrolidine analog of phencyclidine: Some trade or
28 other names: 1-(1-phenylcyclohexyl)pyrrolidine; PCPy; PHP;

29 ~~((+33))~~ (34) Thiophene analog of phencyclidine: Some trade or
30 other names: 1-(1-[2-thienyl]-cyclohexyl)-piperidine; 2-thienyl analog of
31 phencyclidine; TPCP; TCP;

32 ~~((+34))~~ (35) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine: A trade or
33 other name is TCPy.

34 (d) Depressants. Unless specifically excepted or unless listed in
35 another schedule, any material, compound, mixture, or preparation which
36 contains any quantity of the following substances having a depressant
37 effect on the central nervous system, including its salts, isomers, and

1 salts of isomers whenever the existence of such salts, isomers, and
2 salts of isomers is possible within the specific chemical designation.

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

6 (2) Mecloqualone;

7 (3) Methaqualone.

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

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

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

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

18 (4) Fenethylamine;

19 (5) Methcathinone: Some other names: 2-(methylamino)-
20 propiophenone; alpha-(methylamino)propiophenone; 2-(methylamino)-1-
21 phenylpropan-1-one; alpha-N-methylaminopropiophenone;
22 monomethylpropion; ephedrone; N-methylcathinone; methylcathinone; AL-
23 464; AL-422; AL-463 and UR1432, its salts, optical isomers, and salts
24 of optical isomers;

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

27 (7) N-ethylamphetamine;

28 (8) N,N-dimethylamphetamine: Some trade or other names: N,N-
29 alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenoethylene;

30 (9) Any compound structurally derived from 2-amino-1-phenyl-1-
31 propanone (cathinone), including its salts, isomers, and salts of
32 isomers, unless specifically excepted, whenever the existence of such
33 salts, isomers, and salts of isomers is possible within the specific
34 chemical designation, in one or more of the following ways:

35 (i) By substitution of the phenyl ring with alkyl, haloalkyl,
36 alkoxy, alkylendioxy, or halide groups, whether or not further
37 substituted in the phenyl ring to any extent;

38 (ii) By substitution at the 3-position with an alkyl group;

1 (iii) By substitution at the nitrogen atom with alkyl or dialky
2 groups, or by inclusion of the nitrogen in a pyrrolidino or phthalimido
3 group;

4 (iv) The following compounds excepted: Bupropion.

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

7 NEW SECTION. **Sec. 3.** This act is necessary for the immediate
8 preservation of the public peace, health, or safety, or support of the
9 state government and its existing public institutions, and takes effect
10 immediately.

--- END ---