

SENATE JUDICIARY COMMITTEE SUBSTITUTE FOR
SENATE PUBLIC AFFAIRS COMMITTEE SUBSTITUTE FOR
SENATE BILL 127

51ST LEGISLATURE - STATE OF NEW MEXICO - SECOND SESSION, 2014

AN ACT

RELATING TO CONTROLLED SUBSTANCES; AMENDING THE CONTROLLED
SUBSTANCES ACT TO ADD AND REMOVE CERTAIN SUBSTANCES IN THE LIST
OF SCHEDULE I CONTROLLED SUBSTANCES.

BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF NEW MEXICO:

SECTION 1. Section 30-31-6 NMSA 1978 (being Laws 1972,
Chapter 84, Section 6, as amended) is amended to read:

"30-31-6. SCHEDULE I.--The following controlled
substances are included in Schedule I:

A. any of the following opiates, including their
isomers, esters, ethers, salts, and salts of isomers, esters
and ethers, unless specifically exempted, whenever the
existence of these isomers, esters, ethers and salts is
possible within the specific chemical designation:

(1) acetylmethadol;

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underscored material = new
[bracketed material] = delete

underscoring material = new
~~[bracketed material] = delete~~

- 1 (2) allylprodine;
- 2 (3) alphacetylmethadol;
- 3 (4) alphameprodine;
- 4 (5) alphamethadol;
- 5 (6) alpha-methylfentanyl;
- 6 [~~(6)~~] (7) benzethidine;
- 7 [~~(7)~~] (8) betacetylmethadol;
- 8 [~~(8)~~] (9) betameprodine;
- 9 [~~(9)~~] (10) betamethadol;
- 10 [~~(10)~~] (11) betaprodine;
- 11 [~~(11)~~] (12) clonitazene;
- 12 [~~(12)~~] (13) dextromoramide;
- 13 [~~(13)~~] (14) dextrorphan;
- 14 [~~(14)~~] (15) diampromide;
- 15 [~~(15)~~] (16) diethylthiambutene;
- 16 (17) difenoxin;
- 17 [~~(16)~~] (18) dimenoxadol;
- 18 [~~(17)~~] (19) dimepheptanol;
- 19 [~~(18)~~] (20) dimethylthiambutene;
- 20 [~~(19)~~] (21) dioxaphetyl butyrate;
- 21 [~~(20)~~] (22) dipipanone;
- 22 [~~(21)~~] (23) ethylmethylthiambutene;
- 23 [~~(22)~~] (24) etonitazene;
- 24 [~~(23)~~] (25) etoxeridine;
- 25 [~~(24)~~] (26) furethidine;

~~(25)~~ (27) hydroxypethidine;
~~(26)~~ (28) ketobemidone;
~~(27)~~ (29) levomoramide;
~~(28)~~ (30) levophenacymorphan;
~~(29)~~ (31) morpheridine;
~~(30)~~ (32) noracymethadol;
~~(31)~~ (33) norlevorphanol;
~~(32)~~ (34) normethadone;
~~(33)~~ (35) norpipanone;
~~(34)~~ (36) phenadoxone;
~~(34)~~ (37) phenampromide;
~~(36)~~ (38) phenomorphan;
~~(37)~~ (39) phenoperidine;
~~(38)~~ (40) piritramide;
~~(39)~~ (41) proheptazine;
~~(40)~~ (42) properidine;
~~(41)~~ (43) racemoramide; ~~and~~
(44) tilidine; and
~~(42)~~ (45) trimeperidine;

B. any of the following opium derivatives, their salts, isomers and salts of isomers, unless specifically exempted, whenever the existence of these salts, isomers and salts of isomers is possible within the specific chemical designation:

(1) acetorphine;

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underscoring material = new
[bracketed material] = delete

- 1 (2) acetyl dihydrocodeine;
- 2 (3) benzyl morphine;
- 3 (4) codeine methylbromide;
- 4 (5) codeine-N-oxide;
- 5 (6) cyprenorphine;
- 6 (7) desomorphine;
- 7 (8) dihydromorphine;
- 8 (9) etorphine;
- 9 (10) heroin;
- 10 (11) hydromorphinol;
- 11 (12) methyldesorphine;
- 12 (13) methyldihydromorphine;
- 13 (14) morphine methylbromide;
- 14 (15) morphine methylsulfonate;
- 15 (16) morphine-N-oxide;
- 16 (17) myrophine;
- 17 (18) nicocodeine;
- 18 (19) nicomorphine;
- 19 (20) normorphine;
- 20 (21) pholcodine; [~~and~~]
- 21 (22) thebacon;
- 22 (23) drotebanol;
- 23 (24) beta-hydroxy-3-methylfentanyl;
- 24 (25) 3-methylthiofentanyl;
- 25 (26) acetyl-alpha-methylfentanyl;

1 (27) alpha-methylthiofentanyl;

2 (28) beta-hydroxfentanyl;

3 (29) para-fluoro fentanyl; and

4 (30) thiofentanyl;

5 C. any material, compound, mixture or preparation
6 that contains any quantity of the following hallucinogenic
7 substances, their salts, isomers and salts of isomers, unless
8 specifically exempted, whenever the existence of these salts,
9 isomers and salts of isomers is possible within the specific
10 chemical designation:

11 (1) 3,4-methylenedioxy amphetamine;

12 (2) 5-methoxy-3,4-methylenedioxy amphetamine;

13 (3) 3,4,5-trimethoxy amphetamine;

14 (4) bufotenine;

15 (5) diethyltryptamine (DET);

16 (6) dimethyltryptamine (DMT);

17 (7) 4-methyl-2,5-dimethoxy-amphetamine (DOM)

18 or (STP);

19 (8) ibogaine;

20 (9) lysergic acid diethylamide;

21 [~~(10)~~] ~~marijuana;~~

22 [~~(11)~~] (10) mescaline;

23 [~~(12)~~] (11) peyote, except as otherwise

24 provided in the Controlled Substances Act;

25 [~~(13)~~] (12) N-ethyl-3-piperidyl benzilate;

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- 1 [~~(14)~~] (13) N-methyl-3-piperidyl benzilate;
2 [~~(15)~~] (14) psilocybin;
3 [~~(16)~~] (15) psilocyn;
4 [~~(17)~~] (16) tetrahydrocannabinols;
5 [~~(18)~~] (17) hashish;
6 [~~(19) synthetic cannabinoids, including]~~
7 (18) parahexyl (synthetic analog of
8 delta-9-tetra-hydrocannabinol (THC), an active ingredient of
9 cannabis);
10 (19) 2, 5-dimethoxyamphetamine (DMA);
11 (20) 4-bromo-2, 5-dimethoxy-amphetamine (DOB);
12 (21) 4-methoxyamphetamine (PMA);
13 (22) N-ethyl-1-phenylcyclohexylamine (PCE),
14 (ethylamine analog of phencyclidine);
15 (23) 1-(1-phenylcyclohexyl)pyrrolidine (PCPy),
16 (pyrrolidine analog of phencyclidine);
17 (24) thiophene analog of phencyclidine (TCP)
18 or (TPCP);
19 (25) alpha-ethyltryptamine;
20 (26) 2, 5-dimethoxy-4-ethylamphet-amine;
21 (27) ibogaine;
22 (28) 2, 5 dimethoxy-4-(n)-
23 propylthiophenethylamine (2C-T-7);
24 (29) alpha-methyltryptamine (AMT);
25 (30) 5-methoxy-N,N-diisopropyltryptamine

1 (5-MeO-DIPT);

2 (31) synthetic cannabinoids, unless
 3 specifically exempted or unless listed in another schedule,
 4 including any material, compound, mixture or preparation that
 5 contains any quantity of synthetic cannabinoids that
 6 demonstrate binding activity to the cannabinoid receptor or
 7 analogs or homologs with binding activity, including:

8 (a) CP 55,244 ((hydroxymethyl)-4-[2-
 9 hydroxy-4-(2-methyloctan-2-yl)phenyl] 1,2,3,4,4a,5,6,7,8,8a-
 10 decahydronaphthalen-2-ol);

11 (b) CP 55,940 ((5-hydroxy-2-(3-
 12 hydroxypropyl) cyclohexyl)-5-(2-methyloctan-2-yl)phenol);

13 (c) JWH-081 (1-pentyl-3-[1-(4-
 14 methoxynaphthoyl)]indole);

15 (d) JWH-122 (1-pentyl-3-(4-methyl-1-
 16 naphthoyl)indole);

17 (e) JWH-133 3-(1,1-dimethylbutyl)-6a,7,
 18 10,10a-tetrahydro -6,6,9-trimethyl-6H dibenzo[b,d]pyran;

19 (f) JWH 203 (1-pentyl-3-(2-
 20 chlorophenylacetyl)indole);

21 (g) JWH 210 4-ethylnaphthalen-1-yl-(1-
 22 pentylindol-3-yl)methanone;

23 (h) AM-694 (1-(5-fluoropentyl)-3-(2-
 24 iodobenzoyl)indole);

25 (i) AM-1221 (1-(N-methylpiperdin)-2-yl)

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1 methyl-2-methyl-3-(1-naphthoyl)-6-nitroindole;

2 (j) AM-2201 (1-(5-fluoropentyl)-3-(1-
3 naphthoyl)indole);

4 (k) RCS-4 or SR-19 (1-pentyl-3-[(4-
5 methoxy)-benzoyl]indole);

6 (l) RCS-8 or SR-18 (1-cyclohexylethyl-3-
7 (2-methoxyphenylacetyl)indole);

8 (m) WIN-49,098 (pravadoline)(4-
9 methoxyphenyl)-[2-methyl-1-(2-morpholin-4-ylethyl)indol-3-yl]
10 methanone;

11 (n) WIN-55,212-2 (2,3-dihydro-5-methyl-
12 3-(4-morpholinylmethyl)pyrrolo-1,4-benzooxazin-6-yl)-1-
13 naphthalenylmethanone;

14 (o) any of the following synthetic
15 cannabinoids, their salts, isomers and salts of isomers, unless
16 specifically excepted, whenever the existence of these salts,
17 isomers and salts of isomers is possible within the specific
18 chemical designation: 1) naphthoylindoles, or any compound
19 containing a 3-(1-naphthoyl) indole structure with substitution
20 at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
21 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
22 piperidinyl) methyl or 2-(4-morpholinyl) ethyl group, whether
23 or not further substituted in the indole ring to any extent,
24 and whether or not substituted in the naphthyl ring to any
25 extent, including, but not limited to, JWH-015, JWH-018,

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1 JWH-019, JWH-073, JWH-081, JWH-122, JWH-200, JWH-210, JWH-398
2 and AM-2201; 2) naphthylmethylindoles, or any compound
3 containing alhindol-3-yl-(1-naphthyl) methane structure with
4 substitution at the nitrogen atom of the indole ring by an
5 alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
6 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl
7 group, whether or not further substituted in the indole ring to
8 any extent, and whether or not substituted in the naphthyl ring
9 to any extent, including, but not limited to, JWH-175, JWH-184
10 and JWH-199; 3) naphthoylpyrroles, or any compound containing a
11 3-(1-naphthoyl) pyrrole structure with substitution at the
12 nitrogen atom of the pyrrole ring by an alkyl, haloalkyl,
13 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
14 piperidinyl) methyl or 2-(4-morpholinyl) ethyl group, whether
15 or not further substituted in the pyrrole ring to any extent,
16 and whether or not substituted in the naphthyl ring to any
17 extent, including, but not limited to, JWH-307; 4)
18 naphthylmethylindenes, or any compound containing a
19 naphthylideneindene structure with substitution at the
20 3-position of the indene ring by an alkyl, haloalkyl, alkenyl,
21 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)
22 methyl or 2-(4-morpholinyl) ethyl group, whether or not further
23 substituted in the indene ring to any extent, and whether or
24 not substituted in the naphthyl ring to any extent, including,
25 but not limited to, JWH-176; 5) phenylacetylindoles, or any

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1 compound containing a 3-phenylacetylindole structure with
2 substitution at the nitrogen atom of the indole ring by an
3 alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
4 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl
5 group, whether or not further substituted in the indole ring to
6 any extent, and whether or not substituted in the phenyl ring
7 to any extent, including, but not limited to, JWH-203, JWH-250,
8 JWH-251 and RCS-8; 6) cyclohexylphenols, or any compound
9 containing a 2-(3-hydroxycyclohexyl) phenol structure with
10 substitution at the 5- position of the phenolic ring by an
11 alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
12 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl
13 group, whether or not substituted in the cyclohexyl ring to any
14 extent, including, but not limited to, cannabicyclohexanol (CP
15 47,497 C8 homologue), CP 47,497 and CP 55,490; and 7)
16 benzoylindoles, or any compound containing a 3-(benzoyl) [5]
17 OTS-3833.4 indole structure with substitution at the nitrogen
18 atom of the indole ring by an alkyl, haloalkyl, alkenyl,
19 cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)
20 methyl or 2-(4-morpholinyl) ethyl group, whether or not further
21 substituted in the indole ring to any extent, and whether or
22 not substituted in the phenyl ring to any extent, including,
23 but not limited to, AM-694, pravadoline (WIN 48,098), RCS-4 and
24 AM-1241;

25 (p) UR-144 1-(pentyl-1H-indol-3-yl)(2,2,

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1 3,3-tetramethylcyclopropyl)methanone;

2 (q) XLR11 1-(5-fluoro-pentyl)-1H-indol-
3 3-yl(2,2,3,3-tetramethylcyclopropyl)methanone;

4 (r) AKB48 N-(1-adamantyl)-1-pentyl-1H-
5 indazole-3-carboxamide;

6 ~~(a)~~ (s) 1-[2-4-(morpholinyl)ethyl]
7 -3-(1-naphthoyl)indole;

8 ~~(b)~~ (t) 1-butyl-3-(1-naphthoyl)indole;

9 ~~(c)~~ (u) 1-hexyl-3-(1-naphthoyl)indole;

10 ~~(d)~~ (v) 1-pentyl-3-(1-naphthoyl)
11 indole;

12 ~~(e)~~ (w) 1-pentyl-3-(2-
13 methoxyphenylacetyl) indole;

14 ~~(f)~~ (x) cannabicyclohexanol, CP 47,
15 497 and homologues: 5-(1,1-dimethylheptyl)-2-[(1R,3S)
16 -3-hydroxycyclohexyl]-phenol (CP-47,497); and 5-(1,
17 1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol;

18 ~~(g)~~ (y) (6aR,10aR)-9-(hydroxymethyl)
19 -6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
20 tetrahydrobenzo[c]chromen-1-ol;

21 ~~(h)~~ (z) dexanabinol, (6aS,10aS)
22 -9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)
23 -6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;

24 ~~(i)~~ (aa) 1-pentyl-3-(4-chloro
25 naphthoyl) indole;

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1 [~~(j)~~] (bb) (2-methyl-1-propyl-1H-indol-
2 3-yl)-1-naphthalenyl-methanone; and

3 [~~(k)~~] (cc) 5-(1,1-dimethylheptyl)-2-(3-
4 hydroxy cyclohexyl)-phenol;

5 [~~(20)~~] (32) 3,4-methylenedioxymethcathinone;

6 [~~(21)~~] (33) 3,4-methylenedioxyprovalerone;

7 [~~(22)~~] (34) 4-methylmethcathinone;

8 [~~(23)~~] (35) 4-methoxymethcathinone;

9 [~~(24)~~] (36) 3-fluoromethcathinone; [~~and~~

10 [~~(25)~~] (37) 4-fluoromethcathinone;

11 (38) substances determined by the board to

12 have the pharmacological effect of the substance, the risk to
13 the public health by abuse of the substance and the potential
14 of the substance to produce psychic or physiological dependence
15 liability, including:

16 (a) salvia divinorum; and

17 (b) salvinorin A (methyl (2S,4aR,6aR,7R,
18 9S,10aS,10bR)-9-(acetyloxy)-2-(furan-3-yl)-6a,10b-dimethyl-4,10
19 -dioxododecahydro-2H-benzo[f]isochromene-7-carboxylate);

20 (39) 4-methyl-ethylcathinone (4-MEC);

21 (40) 4-ethyl-methcathinone (4-EMC);

22 (41) 2-ethylamino-1-phenyl-propan-1-one
23 (ethcathinone);

24 (42) 3',4'-methylenedioxyethcathinone
25 (ethylone);

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underscored material = new
[bracketed material] = delete

- 1 (43) beta-keto-N-methyl-3,4-
2 benzodioxolybutanamine (bk-MBDB, butylone);
3 (44) naphthylpyrovalerone (NRG-1, naphyrone);
4 (45) N,N-dimethylcathinone (metamfepramone);
5 (46) alpha-pyrrolidinopropiophenone (alpha-
6 PPP);
7 (47) alpha-pyrrolidinobutiophenone (alpha-
8 PBP);
9 (48) 4'-methoxy-alpha-pyrrolidinopropiophenone
10 (MOPPP);
11 (49) 4'-methyl-alpha-pyrrolidinopropiophenone
12 (MPPP);
13 (50) 3',4'-methylenedioxy-alpha-
14 pyrrolidinopropiophenone (MDPPP);
15 (51) 3',4'-methylenedioxy-alpha-
16 pyrrolidinobutiophenone (MDPBP);
17 (52) 4'-methyl-alpha-pyrrolidinobutiophenone
18 (MPBP);
19 (53) alpha-pyrrolidinovalerophenone (alpha-
20 PVP);
21 (54) 5,6-methylenedioxy-2-aminoindane (MDAI);
22 (55) alpha-methylamino-butyrophenone
23 (buphedrone);
24 (56) beta-keto-ethylbenzodioxolybutanamine
25 (eutylone); and

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1 (57) beta-keto-ethylbenzodioxolylpentanamine
2 (pentylone);

3 D. any of the following, unless specifically exempt
4 or unless listed in another schedule, including any material,
5 compound, mixture or preparation that contains any quantity of
6 the following substances having a depressant effect on the
7 central nervous system, including its salts, isomers and salts
8 of isomers, whenever the existence of such salts, isomers and
9 salts of isomers is possible within the specific chemical
10 designation:

11 (1) mecloqualone;

12 (2) methaqualone;

13 (3) benzodiazepines, including:

14 (a) bromazepam;

15 (b) camazepam;

16 (c) cloxazolam;

17 (d) delorazepam;

18 (e) ethylloflazepate;

19 (f) fludiazepam;

20 (g) flunitrazepam;

21 (h) haloxazolam;

22 (i) ketazolam;

23 (j) lopraxolam;

24 (k) lormetazepam;

25 (l) medazepam;

underscored material = new
[bracketed material] = delete

- 1 (m) nimetazepam;
- 2 (n) nitrazepam;
- 3 (o) nordiazepam;
- 4 (p) oxazolam;
- 5 (q) pinazepam; and
- 6 (r) tetrazepam;

7 (4) gamma hydroxybutyric acid and any chemical
8 compound that is metabolically converted to GHB;

9 (5) gamma butyrolactone and any chemical
10 compound that is metabolically converted to GHB; and

11 (6) 1-4 butane diol and any chemical compound
12 that is metabolically converted to GHB;

13 E. any of the following, unless specifically
14 exempted or unless listed in another schedule, including any
15 material, compound, mixture or preparation that contains any
16 quantity of the following substances having a stimulant effect
17 on the central nervous system, including its salts, isomers and
18 salts of isomers:

- 19 (1) fenethylamine;
- 20 (2) N-ethylamphetamine;
- 21 (3) cis-4-methylaminorex;
- 22 (4) N, N-dimethylamphetamine; and
- 23 (5) N-benzylpiperazine (BZP, 1-
24 benzylpiperazine);

25 F. any material, compound, mixture or preparation

1 that contains any quantity of the following substances:

2 (1) 3-methylfentanyl(N-3-methyl-1-(2-phenyl-
3 ethyl)-4-piperidyl)-N-phenylpropanamide, including its optical
4 and geometric isomers, salts and salts of isomers;

5 (2) 3, 4-methylenedioxymethamphetamine (MDMA),
6 including its optical, positional and geometric isomers, salts
7 and salts of isomers;

8 (3) 1-methyl-4-phenyl-4-propionoxypiperidine
9 (MPPP), including its optical isomers, salts, and salts of
10 isomers;

11 (4) 1-(-2-phenylethyl)-4-phenyl-4-acetoxy
12 piperidine (PEPAP), including its optical isomers, salts and
13 salts of isomers;

14 (5) cathinone; and

15 (6) methcathinone;

16 [~~D.~~] G. the enumeration of peyote as a controlled
17 substance does not apply to the use of peyote in bona fide
18 religious ceremonies by a bona fide religious organization, and
19 members of the organization so using peyote are exempt from
20 registration. Any person who manufactures peyote for or
21 distributes peyote to the organization or its members shall
22 comply with the federal Comprehensive Drug Abuse Prevention and
23 Control Act of 1970 and all other requirements of law;

24 [~~E.~~] H. the enumeration of [~~marijuana~~]
25 tetrahydrocannabinols or chemical derivatives of

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1 tetrahydrocannabinol as Schedule I controlled substances does
2 not apply to the use of [~~marijuana~~] tetrahydrocannabinols or
3 chemical derivatives of tetrahydrocannabinol by certified
4 patients pursuant to the Controlled Substances Therapeutic
5 Research Act or by qualified patients pursuant to the
6 provisions of the Lynn and Erin Compassionate Use Act; and
7 [~~F.~~] I. controlled substances added to Schedule I
8 by rule adopted by the board pursuant to Section 30-31-3 NMSA
9 1978."