## SENATE BILL 127

## 51st legislature - STATE OF NEW MEXICO - second session, 2014

## INTRODUCED BY

Sue Wilson Beffort

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AN ACT

RELATING TO CONTROLLED SUBSTANCES; AMENDING THE CONTROLLED SUBSTANCES ACT TO ADD MORE SYNTHETIC CANNABINOIDS AND OTHER SUBSTANCES TO 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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1	(2) allylprodine;
2	(3) alphacetylmethadol;
3	(4) alphameprodine;
4	(5) alphamethadol;
5	(6) alph-methylfentanyl;
6	[ <del>(6)</del> ] <u>(7)</u> benzethidine;
7	[ <del>(7)</del> ] <u>(8)</u> betacetylmethadol;
8	[ <del>(8)</del> ] <u>(9)</u> betameprodine;
9	[ <del>(9)</del> ] <u>(10)</u> betamethadol;
10	[ <del>(10)</del> ] <u>(11)</u> betaprodine;
11	[ <del>(11)</del> ] <u>(12)</u> clonitazene;
12	[ <del>(12)</del> ] <u>(13)</u> dextromoramide;
13	[ <del>(13)</del> ] <u>(14)</u> dextrorphan;
14	[ <del>(14)</del> ] <u>(15)</u> diampromide;
15	[ <del>(15)</del> ] <u>(16)</u> diethylthiambutene;
16	(17) difenoxin;
17	[ <del>(16)</del> ] <u>(18)</u> dimenoxadol;
18	[ <del>(17)</del> ] <u>(19)</u> dimepheptanol;
19	[ <del>(18)</del> ] <u>(20)</u> dimethylthiambutene;
20	[ <del>(19)</del> ] <u>(21)</u> dioxaphetyl butyrate;
21	[ <del>(20)</del> ] <u>(22)</u> dipipanone;
22	[ <del>(21)</del> ] <u>(23)</u> ethylmethylthiambutene;
23	[ <del>(22)</del> ] <u>(24)</u> etonitazene;
24	[ <del>(23)</del> ] <u>(25)</u> etoxeridine;
25	[ <del>(24)</del> ] <u>(26)</u> furethidine;

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 $[\frac{(25)}{(27)}]$ 

[<del>(26)</del>] <u>(28)</u>

 $[\frac{(27)}{(29)}]$ 

hydroxypethidine;

ketobemidone;

levomoramide;

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) dehydro morphine;
9	[ <del>(9)</del> ] <u>(10)</u> etorphine;
10	[ <del>(10)</del> ] <u>(11)</u> heroin;
11	[ <del>(11)</del> ] <u>(12)</u> hydromorphinol;
12	[ <del>(12)</del> ] <u>(13)</u> methyldesorphine;
13	[ <del>(13)</del> ] <u>(14)</u> methyldihydromorphine;
14	[ <del>(14)</del> ] <u>(15)</u> morphine methylbromide;
15	$[\frac{(15)}{(16)}]$ morphine methylsulfonate;
16	[ <del>(16)</del> ] <u>(17)</u> morphine-N-oxide;
17	[ <del>(17)</del> ] <u>(18)</u> myrophine;
18	[ <del>(18)</del> ] <u>(19)</u> nicocodeine;
19	[ <del>(19)</del> ] <u>(20)</u> nicomorphine;
20	[ <del>(20)</del> ] <u>(21)</u> normorphine;
21	[ <del>(21)</del> ] <u>(22)</u> pholcodine; [ <del>and</del>
22	<del>(22)</del> ] <u>(23)</u> thebacon;
23	(24) drotebanol;
24	(25) beta-hydroxy-3-methylfentanyl;
25	(26) 3-methylthiofentanyl;

1	(27) acetyl-alpha-methylfentanyl;
2	(28) alpha-methylthiofentanyl;
3	(29) beta-hydroxfentanyl;
4	(30) para-fluoro fentanyl; and
5	(31) thiofentanyl;
6	C. any material, compound, mixture or preparation
7	that contains any quantity of the following hallucinogenic
8	substances, their salts, isomers and salts of isomers, unless
9	specifically exempted, whenever the existence of these salts,
10	isomers and salts of isomers is possible within the specific
11	chemical designation:
12	(1) 3,4-methylenedioxy amphetamine;
13	(2) 5-methoxy-3,4-methylenedioxy amphetamine;
14	(3) 3,4,5-trimethoxy amphetamine;
15	(4) bufotenine;
16	(5) diethyltryptamine (DET);
17	(6) dimethyltryptamine (DMT);
18	(7) 4-methyl-2,5-dimethoxy-amphetamine (DOM)
19	or (STP);
20	(8) ibogaine;
21	(9) lysergic acid diethylamide;
22	(10) marijuana;
23	(11) mescaline;
24	(12) peyote, except as otherwise provided in
25	the Controlled Substances Act;
	.195503.1

1	(13) N-ethyl-3-piperidyl benzilate;
2	(14) N-methyl-3-piperidyl benzilate;
3	(15) psilocybin;
4	(16) psilocyn;
5	(17) tetrahydrocannabinols;
6	(18) hashish;
7	[ <del>(19) synthetic cannabinoids, including</del> ]
8	(19) parahexyl (synthetic analog of
9	delta-9-tetra-hydrocannabinol (THC), an active ingredient of
10	<pre>cannabis);</pre>
11	(20) 2, 5-dimethoxyamphetamine; 2,5-DMA;
12	(21) 4-bromo-2, 5-dimethoxy-amphetamine;
13	2, 5-DMA;
14	(22) 4-methoxyamphetamine; PMA;
15	(23) ethylamine N-ethyl-1-
16	<pre>phenylcyclohexylamine (PCE);</pre>
17	(24) pyrrolidine l-(1-phenylcyclohexyl)-
18	pyrrolidine (PCPy), (PHP) analog of the drug phencyclidine;
19	(25) thiophene (analog of phencyclidine) (TCP)
20	or (TPCP);
21	(26) alpha-ethyltryptamine;
22	(27) 2, 5-dimethoxy-4-ethylamphet-amine;
23	(28) ibogaine;
24	(29) 2,.5 dimethoxy-4-(n)-
25	<pre>propylthiophenethylamine (2C-T-7);</pre>
	.195503.1

1	(30) alpha-methyltryptamine (AMT);
2	(31) 5-methoxy-N,N-diisopropyltryptamine
3	<u>(5-MeO-DIPT);</u>
4	(32) synthetic cannabinoids, unless
5	specifically exempted or unless listed in another schedule,
6	including any material, compound, mixture or preparation that
7	contains any quantity of synthetic cannabinoids that
8	demonstrate binding activity to the cannabinoid receptor or
9	analogs or homologs with binding activity, including:
10	(a) CP 55,244 ((hydroxymethyl)-4-[2-
11	hydroxy-4-(2-methyloctan-2-yl)phenyl] 1,2,3,4,4a,5,6,7,8,8a-
12	decahydronaphthalen-2-ol);
13	(b) CP 55,940 (5-hydroxy-2-(3-
14	<pre>hydroxypropy1) cyclohexy1]-5-(2-methyloctan-2-y1)phenol);</pre>
15	(c) JWH-081 (1-penty1-3-[1-(4-
16	<pre>methoxynaphthoy)]indole);</pre>
17	(d) JWH-122 (1-penty1-3-(4-methy1-1-
18	<pre>naphthoy1)indole);</pre>
19	(e) JWH-133 3-(1,1-dimethylbutyl)-6a,7,
20	10,10a-tetrahydro -6,6,9-trimethyl-6H dibenzo[b,d]pyran;
21	(f) JWH 203 l-pentyl-3-(2-
22	<pre>chlorophenylacetyl)indole);</pre>
23	(g) JWH 210 4-ethylnaphthalen-1-yl-(1-
24	<pre>pentylindol-3-yl)methanone;</pre>
25	(h) AM-694 (1-(5-fluoropenty1)-3-(2-
	.195503.1

1	<pre>iodobenzoy1)indole);</pre>
2	(i) AM-1221 (1-(N-methylpiperdin-2-yl)
3	<pre>methy1-2-methy1-3-(1-naphthoy1)-6-nitroindole;</pre>
4	(j) AM-2201 (1-(5-fluoropenty1)-3-(1-
5	naphthoyl)indole);
6	(k) RCS-4 or SR-19 (1-penty1-3-[(4-
7	<pre>methoxy)-benzoyl]indole);</pre>
8	(1) RCS-8 or SR-18 (1-cyclohexylethyl-3-
9	(2-methoxyphenylacetyl)indole);
10	(m) JWH-210 (1-penty1-3-(4-
11	<pre>ethylnaphthoyl)indole);</pre>
12	(n) WIN-49,098 (pravadoline)(4-
13	<pre>methoxypheny1)-[2-methy1-1-(2-morpholin-4-ylethy1)indo1-3-y1]</pre>
14	methanone;
15	(o) WIN-55,212-2 (2,3-dihydro-5-methyl-
16	3-(4-morpholinylmethyl)pyrrolo-1,4-benzooxazin6-yl)-1-
17	naphthalenylmethanone);
18	(p) any of the following synthetic
19	cannabinoids, their salts, isomers and salts of isomers, unless
20	specifically excepted, whenever the existence of these salts,
21	isomers and salts of isomers is possible within the specific
22	chemical designation: 1) naphthoylindoles, or any compound
23	containing a 3-(1-naphthoy1) indole structure with substitution
24	at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
25	alkenyl, cycloalkylmethyl, cycloalkylethyl,
	.195503.1

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, including, but not limited to, JWH-015, JWH-018,
<u>JWH-019</u> , <u>JWH-073</u> , <u>JWH-081</u> , <u>JWH-122</u> , <u>JWH-200</u> , <u>JWH-210</u> , <u>JWH-398</u>
and AM-2201; 2) naphthylmethylindoles, or any compound
containing alhindol-3-yl-(1-naphthyl) methane structure with
substitution at the nitrogen atom of the indole ring by an
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, including, but not limited to, JWH-175, JWH-184
and JWH-199; 3) naphthoylpyrroles, or any compound containing a
3-(1-naphthoy1) pyrrole structure with substitution at the
nitrogen atom of the pyrrole ring by an 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,
and whether or not substituted in the naphthyl ring to any
extent, including, but not limited to, JWH-307; 4)
naphthylmethylindenes, or any compound containing a
naphthylideneindene structure with substitution at the
3-position of the indene ring by an alkyl, haloalkyl, alkenyl,
cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)

methyl of 2-(4-morphornyl) ethyl group, whether of not lutther
substituted in the indene ring to any extent, and whether or
not substituted in the naphthyl ring to any extent, including,
but not limited to, JWH-176; 5) phenylacetylindoles, or any
compound containing a 3-phenylacetylindole structure with
substitution at the nitrogen atom of the indole ring by an
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, including, but not limited to, JWH-203, JWH-250,
JWH-251 and RCS-8; 6) cyclohexylphenols, or any compound
containing a 2-(3-hydroxycyclohexyl) phenol structure with
substitution at the 5- position of the phenolic ring by an
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, including, but not limited to, cannabicyclohexanol (CP
47,497 C8 homologue), CP 47,497 and CP 55,490; and 7)
<pre>benzoylindoles, or any compound containing a 3-(benzoyl) [ 5 ]</pre>
OTS-3833.4 indole structure with substitution at the nitrogen
atom of the indole ring by an alkyl, haloalkyl, alkenyl,
cycloalkylmethyl, cycloalkylethyl, l-(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
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       not substituted in the phenyl ring to any extent, including,
       but not limited to, AM-694, pravadoline (WIN 48,098), RCS-4 and
 2
 3
       AM-1241;
                                (q) UR-144 l-(pentyl-lH-indol-3-yl)(2,2,
 4
       3,3-tetramethylcyclopropyl)methanone;
 5
                                (r) XLR11 1-(5-fluoro-penty1)-lH-indol-
 6
 7
       3-y1(2,2,3,3-tetramethylcyclopropyl)methanone;
                                (s) AKB48 N-(1-adamanty1)-1-penty1-1H-
 8
       indazole-3-carboxamide;
 9
                                \lceil \frac{a}{a} \rceil (t) 1-\lceil 2 - (4 - (morpholiny1))  ethy1
10
       -3-(1-naphthoy1)indole;
11
12
                                [\frac{(b)}{(u)}] (u) 1-buty1-3-(1-napthoy1)indole;
                                [\frac{(c)}{(v)}] 1-hexy1-3-(1-naphthoy1)indole;
13
14
                                \left[\frac{\text{(d)}}{\text{(w)}}\right] 1-penty1-3-(1-naphthoy1)
       indole;
15
                                [(e)] (x) 1-penty1-3-(2-
16
       methoxyphenylacetyl) indole;
17
                                [\frac{f}{f}] (y) cannabicyclohexanol (CP 47,
18
       497 and homologues: 5-(1,1-dimethylheptyl)-2-[(1R,3S)]
19
20
       -3-hydroxycyclohexyl]-phenol (CP-47,497); and 5-(1,
       1-dimethyloctyl)-2-[(lR,3S)-3-hydroxycyclohexyl]-phenol;
21
                                [\frac{g}{2}] (z) 6aR, 10aR) -9-(hydroxymethy1)
22
       -6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,
23
       10a-tetrahydrobenzo[c]chromen-1-o1);
24
                                [<del>(h)</del>] <u>(aa)</u> dexanabinol, (6aS,10aS)
25
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1
       -9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)
 2
       -6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;
                               [\frac{(i)}{(b)}] 1-penty1-3-(4-chloro
 3
       naphthoyl) indole;
 4
                               [\frac{(i)}{(i)}] (cc) (2-methyl-l-propyl-lH-indol-
 5
       3-y1)-1-naphthalenyl-methanone; and
 6
 7
                               [\frac{(k)}{(dd)}] 5-(1,1-dimethylheptyl)-2-(3-
       hydroxy cyclohexyl)-phenol;
 8
 9
                         [\frac{(20)}{(33)}] 3,4-methylenedioxymethcathinone;
                         \lceil \frac{(21)}{3} \rceil (34) 3,4-methylenedioxypyrovalerone;
10
                         [\frac{(22)}{(35)}] 4-methylmethcathinone;
11
12
                         [<del>(23)</del>] <u>(36)</u> 4-methoxymethcathinone;
                         [<del>(24)</del>] (37) 3-fluoromethcathinone; [and
13
14
                         (25) (38) 4-fluoromethcathinone;
                         (39) substances determined by the board to
15
       have the pharmacological effect of the substance, the risk to
16
       the public health by abuse of the substance and the potential
17
       of the substance to produce psychic or physiological dependence
18
       liability, including:
19
20
                               (a) salvia divinorum; and
                               (b) salvinorin A (methyl (2S, 4aR, 6aR, 7R,
21
       9S, 10aS, 10bR)-9-(acetyloxy)-2-(furan-3-y1)-6a, 10b-dimethyl-4, 10
22
       -dioxododecahydro-2H-benzo[f]isochromene-7-carboxylate);
23
                         (40) 4-methyl-ethylcathinone (4-MEC);
24
                         (41) 4-ethyl-methcathinone (4-EMC);
25
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1	(42) 2-ethylamino-l-phenyl-propan-l-one
2	<pre>(ethcathinone);</pre>
3	(43) 3',4'-methylenedioxyethcathinone
4	<pre>(ethylone);</pre>
5	(44) beta-keto-N-methyl-3,4-
6	benzodioxyolybutanamine (bk-MBDB, butylone);
7	(45) naphthylpyrovalerone (NRG-1, naphyrone);
8	(46) N,N-dimethylcathinone (metamfepramone);
9	(47) alpha-pyrrolidinopropiophenone
10	(alpha-PPP);
11	(48) alpha-pyrrolidinobutiophenone (?-PBP);
12	(49) 4'-methoxy-alpha-pyrrolidinopropiophenone
13	(MOPPP);
14	(50) 4'-methyl-?-pyrrolidinopropiophenone
15	(MPPP);
16	(51) 3',4'-methylenedioxy-alpha-
17	<pre>pyrrolidinopropiophenone (MDPPP);</pre>
18	(52) 3',4'-methylenedioxy-alpha-
19	<pre>pyrrolidinobutiophenone (MDPBP);</pre>
20	(53) 4'-methyl-?-pyrrolidinobutiophenone
21	(MPBP);
22	(54) alpha-pyrrolidinovalerophenone
23	(alpha-PVP);
24	(55) 5,6-methylenedioxy-2-aminoindane (MDAI);
25	(56) alpha-methylamino-butyrophenone
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1	(buphedrone);
2	(57) beta-keto-ethylbenzodioxolylbutanamine
3	(eutylone); and
4	(58) beta-keto-ethylbenzodioxolylpentanamine
5	<pre>(pentylone);</pre>
6	D. any of the following, unless specifically exempt
7	or unless listed in another schedule, including any material,
8	compound, mixture or preparation that contains any quantity of
9	the following substances having a depressant effect on the
10	central nervous system, including its salts, isomers and salts
11	of isomers, whenever the existence of such salts, isomers and
12	salts of isomers is possible within the specific chemical
13	designation:
14	(1) mecloqualone;
15	(2) methaqualone;
16	(3) benzodiazepines, including:
17	(a) bromazepam;
18	(b) camazepam;
19	(c) cloxazolam;
20	(d) delorazepam;
21	(e) ethylloflazepate;
22	(f) fludiazepam;
23	(g) flunitrazepam;
24	(h) haloxazolam;
25	(i) ketazolam;

1	(J) loprazolam;
2	(k) lormetazepam;
3	(1) medazepam;
4	(m) nimetazepam;
5	(n) nitrazepam;
6	(o) nordiazepam;
7	(p) oxazolam;
8	(q) pinazepam; and
9	(r) tetrazepam;
10	(4) gamma hydroxybutyric acid and any chemical
11	compound that is metabolically converted to GHB;
12	(5) gamma butyrolactone and any chemical
13	compound that is metabolically converted to GHB; and
14	(6) 1-4 butane diol and any chemical compound
15	that is metabolically converted to GHB;
16	E. any of the following, unless specifically
17	exempted or unless listed in another schedule, including any
18	material, compound, mixture or preparation that contains any
19	quantity of the following substances having a stimulant effect
20	on the central nervous system, including its salts, isomers and
21	salts of isomers:
22	(1) fenethylline;
23	(2) N-ethylamphetamine;
24	(3) cis-4-methylaminorex;
25	(4) N, N-dimethylamphetamine; and
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1	(5) N-benzylpiperazine (BZP, 1-							
2	<pre>benzylpiperazine);</pre>							
3	F. any material, compound, mixture or preparation							
4	that contains any quantity of the following substances:							
5	(1) 3-methylfentanyl(N-3-methyl-1-(2-phenyl-							
6	ethyl)-4-piperidyl)-N-phenylpropanamide, including its optical							
7	and geometric isomers, salts and salts of isomers;							
8	(2) 3, 4-methylenedioxymethamphetamine (MDMA),							
9	including its optical, positional and geometric isomers, salts							
10	and salts of isomers;							
11	(3) 1-methyl-4-phenyl-4-proprionoxypiperidine							
12	(MPPP), including its optical isomers, salts, and salts of							
13	<pre>isomers;</pre>							
14	(4) 1-(-2-phenylethyl)-4-phenyl-4-acetoxy							
15	piperidine (PEPAP), including its optical isomers, salts and							
16	salts of isomers;							
17	(5) cathinone; and							
18	(6) methcathinone;							
19	$[\frac{D_{\bullet}}{G_{\bullet}}]$ the enumeration of peyote as a controlled							
20	substance does not apply to the use of peyote in bona fide							
21	religious ceremonies by a bona fide religious organization, and							
22	members of the organization so using peyote are exempt from							
23	registration. Any person who manufactures peyote for or							
24	distributes peyote to the organization or its members shall							
25	comply with the federal Comprehensive Drug Abuse Prevention and							

Control	Act	of	1970	and	a11	other	requirements	of	law;
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[E.] H. the enumeration of marijuana, tetrahydrocannabinols or chemical derivatives of tetrahydrocannabinol as Schedule I controlled substances does not apply to the use of marijuana, tetrahydrocannabinols or chemical derivatives of tetrahydrocannabinol by certified patients pursuant to the Controlled Substances Therapeutic Research Act or by qualified patients pursuant to the provisions of the Lynn and Erin Compassionate Use Act; and

 $[F_{\bullet}]$  I. controlled substances added to Schedule I by rule adopted by the board pursuant to Section 30-31-3 NMSA 1978."

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