1	SENATE PUBLIC AFFAIRS COMMITTEE SUBSTITUTE FOR SENATE BILL 127
2	51st legislature - STATE OF NEW MEXICO - second session, 2014
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10	AN ACT
11	RELATING TO CONTROLLED SUBSTANCES; AMENDING THE CONTROLLED
12	SUBSTANCES ACT TO ADD MORE SYNTHETIC CANNABINOIDS AND OTHER
13	SUBSTANCES TO THE LIST OF SCHEDULE I CONTROLLED SUBSTANCES.
14	
15	BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF NEW MEXICO:
16	SECTION 1. Section 30-31-6 NMSA 1978 (being Laws 1972,
17	Chapter 84, Section 6, as amended) is amended to read:
18	"30-31-6. SCHEDULE IThe following controlled
19	substances are included in Schedule I:
20	A. any of the following opiates, including their
21	isomers, esters, ethers, salts, and salts of isomers, esters
22	and ethers, unless specifically exempted, whenever the
23	existence of these isomers, esters, ethers and salts is
24	possible within the specific chemical designation:
25	<pre>(1) acetylmethadol;</pre>
	.196388.2

[<del>bracketed material</del>] = delete <u>underscored material = new</u>

1	(2) allylprodine;
2	(3) alphacetylmethadol;
3	(4) alphameprodine;
4	(5) alphamethadol;
5	(6) alpha-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	<u>(17) difenoxin;</u>
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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;

	1	[ <del>(25)</del> ] <u>(27)</u> hydroxypethidine;
	2	[ <del>(26)</del> ] <u>(28)</u> ketobemidone;
	3	[ <del>(27)</del> ] <u>(29)</u> levomoramide;
	4	[ <del>(28)</del> ] <u>(30)</u> levophenacylmorphan;
	5	[ <del>(29)</del> ] <u>(31)</u> morpheridine;
	6	[ <del>(30)</del> ] <u>(32)</u> noracymethadol;
	7	[ <del>(31)</del> ] <u>(33)</u> norlevorphanol;
	8	[ <del>(32)</del> ] <u>(34)</u> normethadone;
	9	[ <del>(33)</del> ] <u>(35)</u> norpipanone;
	10	[ <del>(34)</del> ] <u>(36)</u> phenadoxone;
	11	[ <del>(34)</del> ] <u>(37)</u> phenampromide;
	12	[ <del>(36)</del> ] <u>(38)</u> phenomorphan;
	13	[ <del>(37)</del> ] <u>(39)</u> phenoperidine;
	14	[ <del>(38)</del> ] <u>(40)</u> piritramide;
	15	[ <del>(39)</del> ] <u>(41)</u> proheptazine;
	16	[ <del>(40)</del> ] <u>(42)</u> properidine;
ברם	17	[ <del>(41)</del> ] <u>(43)</u> racemoramide; [ <del>and</del> ]
רטח	18	(44) tilidine; and
	19	[ <del>(42)</del> ] <u>(45)</u> trimeperidine;
	20	B. any of the following opium derivatives, their
ווומ רב	21	salts, isomers and salts of isomers, unless specifically
	22	exempted, whenever the existence of these salts, isomers and
l ntacvered	23	salts of isomers is possible within the specific chemical
ן טו מ	24	designation:
	25	(1) acetorphine;
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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; [ <del>and</del> ]
21	(22)	thebacon;
22	<u>(23)</u>	drotebanol;
23	<u>(24)</u>	<pre>beta-hydroxy-3-methylfentanyl;</pre>
24	<u>(25)</u>	3-methylthiofentanyl;
25	<u>(26)</u>	<pre>acety1-alpha-methylfentanyl;</pre>

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1	(27) alpha-methylthiofentanyl;
2	<u>(28) beta-hydroxfentanyl;</u>
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	<ol> <li>3,4-methylenedioxy amphetamine;</li> </ol>
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	<u>or (STP)</u> ;
19	<pre>(8) ibogaine;</pre>
20	(9) lysergic acid diethylamide;
21	(10) marijuana;
22	(11) mescaline;
23	(12) peyote, except as otherwise provided in
24	the Controlled Substances Act;
25	(13) N-ethyl-3-piperidyl benzilate;
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1	(14) N-methyl-3-piperidyl benzilate;
2	(15) psilocybin;
3	<pre>(16) psilocyn;</pre>
4	(17) tetrahydrocannabinols;
5	(18) hashish;
6	[ <del>(19) synthetic cannabinoids, including</del> ]
7	(19) parahexyl (synthetic analog of
8	delta-9-tetra-hydrocannabinol (THC), an active ingredient of
9	<pre>cannabis);</pre>
10	(20) 2, 5-dimethoxyamphetamine (DMA);
11	(21) 4-bromo-2, 5-dimethoxy-amphetamine (DOB);
12	(22) 4-methoxyamphetamine (PMA);
13	(23) N-ethyl-l-phenylcyclohexylamine (PCE),
14	(ethylamine analog of phencyclidine);
15	(24) 1-(1-phenylcyclohexyl)pyrrolidine (PCPy),
16	(pyrrolidine analog of phencyclidine);
17	(25) thiophene analog of phencyclidine (TCP)
18	<u>or (TPCP);</u>
19	(26) alpha-ethyltryptamine;
20	(27) 2, 5-dimethoxy-4-ethylamphet-amine;
21	(28) ibogaine;
	(29) 2, 5 dimethoxy-4-(n)-
22	
22 23	propylthiophenethylamine (2C-T-7);
	propylthiophenethylamine (2C-T-7); (30) alpha-methyltryptamine (AMT);
23	

<u>underscored material = new</u> [<del>bracketed material</del>] = delete

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	1	<u>(5-MeO-DIPT);</u>		
	2	(32) 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	<u>(a) CP 55,244 ((hydroxymethyl)-4-[2-</u>		
	9	hydroxy-4-(2-methyloctan-2-yl)phenyl] 1,2,3,4,4a,5,6,7,8,8a-		
	10	<pre>decahydronaphthalen-2-ol);</pre>		
	11	<u>(b)</u> CP 55,940 ((5-hydroxy-2-(3-		
	12	<pre>hydroxypropy1) cyclohexy1)-5-(2-methyloctan-2-y1)phenol);</pre>		
	13	<u>(c) JWH-081 (l-penty1-3-[l-(4-</u>		
	14	<pre>methoxynaphthoy1)]indole);</pre>		
	15	(d) JWH-122 (1-penty1-3-(4-methy1-1-		
	16	<pre>naphthoyl)indole);</pre>		
	17	<u>(e) JWH-133 3-(1,1-dimethylbutyl)-6a,7,</u>		
	18	<pre>10,10a-tetrahydro -6,6,9-trimethyl-6H dibenzo[b,d]pyran;</pre>		
ı	19	<u>(f) JWH 203 (1-penty1-3-(2-</u>		
	20	<pre>chlorophenylacetyl)indole);</pre>		
	21	<u>(g) JWH 210 4-ethylnaphthalen-1-yl-(1-</u>		
	22	<pre>pentylindol-3-yl)methanone;</pre>		
	23	<u>(h) AM-694 (l-(5-fluoropentyl)-3-(2-</u>		
I	24	<pre>iodobenzoy1)indole);</pre>		
	25	<u>(i) AM-1221 (l-(N-methylpiperdin)-2-yl)</u>		
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## [<del>bracketed material</del>] = delete <u>underscored material = new</u>

1	<pre>methy1-2-methy1-3-(1-naphthoy1)-6-nitroindole;</pre>
2	<u>(j) AM-2201 (1-(5-fluoropentyl)-3-(1-</u>
3	<pre>naphthoyl)indole);</pre>
4	<u>(k) RCS-4 or SR-19 (1-penty1-3-[(4-</u>
5	<pre>methoxy)-benzoyl]indole);</pre>
6	(1) RCS-8 or SR-18 (1-cyclohexylethyl-3-
7	<pre>(2-methoxyphenylacetyl)indole);</pre>
8	(m) WIN-49,098 (pravadoline)(4-
9	<pre>methoxyphenyl)-[2-methyl-l-(2-morpholin-4-ylethyl)indol-3-yl]</pre>
10	<u>methanone;</u>
11	<u>(n) WIN-55,212-2 (2,3-dihydro-5-methyl-</u>
12	<u>3-(4-morpholinylmethyl)pyrrolo-1,4-benzooxazin6-yl)-1-</u>
13	naphthalenylmethanone;
14	(o) any of the following synthetic
14 15	(o) any of the following synthetic cannabinoids, their salts, isomers and salts of isomers, unless
15	cannabinoids, their salts, isomers and salts of isomers, unless
15 16	cannabinoids, their salts, isomers and salts of isomers, unless specifically excepted, whenever the existence of these salts,
15 16 17	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
15 16 17 18	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) naphthoylindoles, or any compound
15 16 17 18 19	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) naphthoylindoles, or any compound containing a 3-(1-naphthoyl) indole structure with substitution
15 16 17 18 19 20	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) naphthoylindoles, or any compound containing a 3-(1-naphthoyl) indole structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
15 16 17 18 19 20 21	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) naphthoylindoles, or any compound containing a 3-(1-naphthoyl) indole structure with substitution at the nitrogen atom of the indole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
15 16 17 18 19 20 21 21 22	<pre>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) naphthoylindoles, or any compound containing a 3-(1-naphthoyl) indole 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</pre>
15 16 17 18 19 20 21 22 23	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) naphthoylindoles, or any compound containing a 3-(1-naphthoyl) indole 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,

[<del>bracketed material</del>] = delete

<u>underscored material = new</u>

	1	<u>JWH-019, JWH-073, JWH-081, JWH-122, JWH-200, JWH-210, JWH-398</u>
	2	and AM-2201; 2) naphthylmethylindoles, or any compound
	3	containing alhindol-3-yl-(l-naphthyl) methane structure with
	4	substitution at the nitrogen atom of the indole ring by an
	5	<u>alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,</u>
	6	<u>l-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl</u>
	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	<u>3-(1-naphthoyl) pyrrole structure with substitution at the</u>
	12	nitrogen atom of the pyrrole ring by an alkyl, haloalkyl,
	13	<u>alkenyl, cycloalkylmethyl, cycloalkylethyl, l-(N-methyl-2-</u>
	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
ר ר ר	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	<u>3-position of the indene ring by an alkyl, haloalkyl, alkenyl, </u>
	21	<u>cycloalkylmethyl, cycloalkylethyl, l-(N-methyl-2-piperidinyl)</u>
- - -	22	methyl or 2-(4-morpholinyl) ethyl group, whether or not further
	23	substituted in the indene ring to any extent, and whether or
5	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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<u>underscored material = new</u> [bracketed material] = delete

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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	<u>alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,</u>
4	<u>l-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl</u>
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	<u>alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,</u>
12	<u>l-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl) ethyl</u>
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, l-(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	<u>AM-1241;</u>
25	<u>(p) UR-144 l-(pentyl-1H-indol-3-yl)(2,2</u>

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1	3,3-tetramethylcyclopropyl)methanone;
2	<u>(q) XLR11 l-(5-fluoro-pentyl)-lH-indol-</u>
3	<u>3-y1(2,2,3,3-tetramethylcyclopropyl)methanone;</u>
4	<u>(r) AKB48 N-(l-adamantyl)-l-pentyl-lH-</u>
5	<u>indazole-3-carboxamide;</u>
6	[ <del>(a)</del> ] <u>(s)</u> l-[2-4-(morpholinyl)ethyl]
7	-3-(l-naphthoyl)indole;
8	<pre>[(b)] (t) l-butyl-3-(l-napthoyl)indole;</pre>
9	<pre>[(c)] (u) l-hexyl-3-(l-naphthoyl)indole;</pre>
10	[ <del>(d)</del> ] <u>(v)</u> l-pentyl-3-(l-naphthoyl)
11	indole;
12	[ <del>(e)</del> ] <u>(w)</u> l-pentyl-3-(2-
13	<pre>methoxyphenylacetyl) indole;</pre>
14	[ <del>(f)</del> ] <u>(x)</u> 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	<pre>l-dimethyloctyl)-2-[(lR,3S)-3-hydroxycyclohexyl]-phenol;</pre>
18	[ <del>(g)</del> ] <u>(y)</u> (6aR,10aR)-9-(hydroxymethy1)
19	-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-
20	<pre>tetrahydrobenzo[c]chromen-l-o1;</pre>
21	[ <del>(h)</del> ] <u>(z)</u> dexanabinol, (6aS,10aS)
22	-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)
23	-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-o1;
24	[ <del>(i)</del> ] <u>(aa)</u> l-pentyl-3-(4-chloro
25	naphthoyl) indole;
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[bracketed material] = delete

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1	
1	[ <del>(j)</del> ] <u>(bb)</u> (2-methyl-l-propyl-lH-indol-
2	3-y1)-1-naphthaleny1-methanone; and
3	[ <del>(k)</del> ] <u>(cc)</u> 5-(1,1-dimethylheptyl)-2-(3-
4	hydroxy cyclohexyl)-phenol;
5	[ <del>(20)</del> ] <u>(33)</u> 3,4-methylenedioxymethcathinone;
6	[ <del>(21)</del> ] <u>(34)</u> 3,4-methylenedioxypyrovalerone;
7	[ <del>(22)</del> ] <u>(35)</u> 4-methylmethcathinone;
8	[ <del>(23)</del> ] <u>(36)</u> 4-methoxymethcathinone;
9	[ <del>(24)</del> ] <u>(37)</u> 3-fluoromethcathinone; [ <del>and</del>
10	(25)] (38) 4-fluoromethcathinone;
11	(39) 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	<u>liability, including:</u>
16	(a) salvia divinorum; and
17	<u>(b) salvinorin A (methyl (2S,4aR,6aR,7R,</u>
18	<u>98,10aS,10bR)-9-(acetyloxy)-2-(furan-3-y1)-6a,10b-dimethyl-4,10</u>
19	<pre>-dioxododecahydro-2H-benzo[f]isochromene-7-carboxylate);</pre>
20	(40) 4-methyl-ethylcathinone (4-MEC);
21	(41) 4-ethyl-methcathinone (4-EMC);
22	<u>(42) 2-ethylamino-l-phenyl-propan-l-one</u>
23	<pre>(ethcathinone);</pre>
24	(43) 3',4'-methylenedioxyethcathinone
25	<pre>(ethylone);</pre>
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1	(44) beta-keto-N-methyl-3,4-
2	<pre>benzodioxyolybutanamine (bk-MBDB, butylone);</pre>
3	(45) naphthylpyrovalerone (NRG-1, naphyrone);
4	(46) N,N-dimethylcathinone (metamfepramone);
5	<u>(47) alpha-pyrrolidinopropiophenone (alpha-</u>
6	<u>PPP);</u>
7	<u>(48) alpha-pyrrolidinobutiophenone (alpha-</u>
8	<u>PBP);</u>
9	(49) 4'-methoxy-alpha-pyrrolidinopropiophenone
10	<u>(MOPPP);</u>
11	(50) 4'-methyl-alpha-pyrrolidinopropiophenone
12	<u>(MPPP);</u>
13	(51) 3',4'-methylenedioxy-alpha-
14	pyrrolidinopropiophenone (MDPPP);
15	(52) 3',4'-methylenedioxy-alpha-
16	pyrrolidinobutiophenone (MDPBP);
17	(53) 4'-methyl-alpha-pyrrolidinobutiophenone
18	<u>(MPBP);</u>
19	<u>(54) alpha-pyrrolidinovalerophenone (alpha-</u>
20	<u>PVP);</u>
21	(55) 5,6-methylenedioxy-2-aminoindane (MDAI);
22	(56) alpha-methylamino-butyrophenone
23	(buphedrone);
24	(57) beta-keto-ethylbenzodioxolylbutanamine
25	(eutylone); and
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1	(58) beta-keto-ethylbenzodioxolylpentanamine
2	<u>(pentylone);</u>
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	<u>(a) bromazepam;</u>
15	<u>(b) camazepam;</u>
16	(c) cloxazolam;
17	(d) delorazepam;
18	(e) ethylloflazepate;
19	<u>(f) fludiazepam;</u>
20	(g) flunitrazepam;
21	(h) haloxazolam;
22	<u>(i) ketazolam;</u>
23	<u>(j) loprazolam;</u>
24	<u>(k) lormetazepam;</u>
25	<u>(1) medazepam;</u>

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1	(m) nimetazepam;
2	<u>(n) nitrazepam;</u>
3	<u>(o) nordiazepam;</u>
4	<u>(p) oxazolam;</u>
5	(q) pinazepam; and
6	<u>(r) tetrazepam;</u>
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	<u>salts of isomers:</u>
19	(1) fenethylline;
20	(2) N-ethylamphetamine;
21	(3) cis-4-methylaminorex;
22	(4) N, N-dimethylamphetamine; and
23	(5) N-benzylpiperazine (BZP, 1-
24	<pre>benzylpiperazine);</pre>
25	F. any material, compound, mixture or preparation
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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) l-methyl-4-phenyl-4-proprionoxypiperidine
9	(MPPP), including its optical isomers, salts, and salts of
10	<u>isomers;</u>
11	(4) l-(-2-phenylethyl)-4-phenyl-4-acetoxy
12	piperidine (PEPAP), including its optical isomers, salts and
13	<u>salts of isomers;</u>
14	(5) cathinone; and
15	(6) methcathinone;
16	$[D_{\cdot}]$ <u>G.</u> 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	[ <del>E.</del> ] <u>H.</u> the enumeration of marijuana,
25	tetrahydrocannabinols or chemical derivatives of

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underscored material = new
[bracketed material] = delete

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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	$[F_{\bullet}]$ <u>I.</u> controlled substances added to Schedule I
8	by rule adopted by the board pursuant to Section 30-31-3 NMSA
9	1978."
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