1	STATE OF OKLAHOMA
2	1st Session of the 60th Legislature (2025)
3	HOUSE BILL 1311 By: Humphrey
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6	AS INTRODUCED
7	An Act relating to controlled dangerous substances;
8	amending 63 O.S. 2021, Section 2-204, as last amended by Section 3, Chapter 308, O.S.L. 2024 (63 O.S. Supp.
9	2024, Section 2-204), which relates to the Uniform Controlled Dangerous Substances Act; adding chemicals
10	to Schedule I; and providing an effective date.
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13	BE IT ENACTED BY THE PEOPLE OF THE STATE OF OKLAHOMA:
14	SECTION 1. AMENDATORY 63 O.S. 2021, Section 2-204, as
15	last amended by Section 3, Chapter 308, O.S.L. 2024 (63 O.S. Supp.
16	2024, Section 2-204), is amended to read as follows:
17	Section 2-204. The controlled substances listed in this section
18	are included in Schedule I and include any material, compound,
19	mixture or preparation that contains any quantity of the following
20	hallucinogenic substances, their salts, isomers and salts of
21	isomers, unless specifically excepted, when the existence of these
22	salts, isomers and salts of isomers is possible within the specific
23	chemical designation.

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

- 6 1. Acetylmethadol;
- 7 2. Allylprodine;
- 8 3. Alphacetylmethadol;
- 9 4. Alphameprodine;
- 10 5. Alphamethadol;
- 11 6. Benzethidine;
- 12 7. Betacetylmethadol;
- 13 8. Betameprodine;
- 14 9. Betamethadol;
- 15 10. Betaprodine;
- 16 11. Clonitazene;
- 17 12. Dextromoramide;
- 18 13. Dextrorphan (except its methyl ether);
- 19 14. Diampromide;
- 20 15. Diethylthiambutene;
- 21 16. Dimenoxadol;
- 22 17. Dimepheptanol;
- 23 18. Dimethylthiambutene;
- 24 19. Dioxaphetyl butyrate;

1	20.	Dipipanone;
2	21.	Ethylmethylthiambutene;
3	22.	Etonitazene;
4	23.	Etoxeridine;
5	24.	Furethidine;
6	25.	Hydroxypethidine;
7	26.	Isotonitazene;
8	27.	Ketobemidone;
9	28.	Levomoramide;
10	29.	Levophenacylmorphan;
11	30.	Metonitazene;
12	31.	Morpheridine;
13	32.	N-desethyl isotonitazene;
14	33.	N-pyrrolidino protonitazene;
15	34.	Noracymethadol;
16	35.	Norlevorphanol;
17	36.	Normethadone;
18	37.	Norpipanone;
19	38.	Phenadoxone;
20	39.	Phenampromide;
21	40.	Phenomorphan;
22	41.	Phenoperidine;
23	42.	Piritramide;
24	43.	Proheptazine;

1	44. Properidine;
2	45. Protonitazene;
3	46. Racemoramide; or
4	47. Trimeperidine.
5	B. Any of the following opium derivatives, their salts,
6	isomers, and salts of isomers, unless specifically excepted, when
7	the existence of these salts, isomers, and salts of isomers is
8	possible within the specific chemical designation:
9	1. Acetorphine;
10	2. Acetyldihydrocodeine;
11	3. Benzylmorphine;
12	4. Codeine methylbromide;
13	5. Codeine-N-Oxide;
14	6. Cyprenorphine;
15	7. Desomorphine;
16	8. Dihydromorphine;
17	9. Etorphine;
18	10. Heroin;
19	11. Hydromorphinol;
20	12. Methyldesorphine;
21	13. Methylhydromorphine;
22	14. Morphine methylbromide;
23	15. Morphine methylsulfonate;
24	16. Morphine-N-Oxide;

1	1 17. Myrophine;	
2	2 18. Nicocodeine;	
3	3 19. Nicomorphine;	
4	4 20. Normorphine;	
5	5 21. Phoclodine;	
6	6 22. Thebacon;	
7	7 23. N-phenyl-N-[1-(2-phenyle	thyl)-4-piperidinyl]-acetamide
8	8 (Acetyl fentanyl);	
9	9 24. N-phenyl-N-[1-(2-phenyle	thyl)-4-piperidinyl]-butenamide
10	0 (Crotonyl fentanyl);	
11	1 25. N-phenyl-N-[1-(2-phenyle	thyl)-4-piperidinyl]-2-
12	2 furancarboxamide (Furanyl fentany	1);
13	3 26. N-phenyl-1-(2-phenylethy	l)-4-piperidinamine (4-ANPP);
14	4 27. N-(1-phenethylpiperidin-	4-yl)-N-
15	5 phenylcyclopropanecarboxamide (Cy	clopropyl fentanyl); or
16	6 28. N-phenyl-N-[1-(2-phenyle	thyl)-4-piperidinyl]-butanamide
17	7 (Butyrl fentanyl).	
18	8 C. Any material, compound, m	ixture, or preparation which
19	9 contains any quantity of the foll	owing hallucinogenic substances,
20	0 their salts, isomers, and salts o	f isomers, unless specifically
21	1 excepted, when the existence of t	hese salts, isomers, and salts of
22	2 isomers is possible within the sp	ecific chemical designation:
23	3 1. Methcathinone;	
24	4 2. 3, 4-methylenedioxy amphe	tamine;

1	3.	3, 4-methylenedioxy methamphetamine;
2	4.	5-methoxy-3, 4-methylenedioxy amphetamine;
3	5.	3, 4, 5-trimethoxy amphetamine;
4	6.	Bufotenine;
5	7.	Diethyltryptamine;
6	8.	Dimethyltryptamine;
7	9.	4-methyl-2, 5-dimethoxyamphetamine;
8	10.	Ibogaine;
9	11.	Lysergic acid diethylamide;
10	12.	Marijuana;
11	13.	Mescaline;
12	14.	N-benzylpiperazine;
13	15.	N-ethyl-3-piperidyl benzilate;
14	16.	N-methyl-3-piperidyl benzilate;
15	17.	Psilocybin;
16	18.	Psilocyn;
17	19.	2, 5 dimethoxyamphetamine;
18	20.	4 Bromo-2, 5-dimethoxyamphetamine;
19	21.	4 methoxyamphetamine;
20	22.	Cyclohexamine;
21	23.	Salvia Divinorum;
22	24.	Salvinorin A;
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1	25.	Thiophene Analog of Phencyclidine. Also known as: 1-(1-(2-					
2	thienyl)	l) cyclohexyl) piperidine; 2-Thienyl Analog of Phencyclidine;					
3	TPCP, TCP;						
4	26.	Phencyclidine (PCP);					
5	27.	Pyrrolidine Analog for Phencyclidine. Also known as 1-(1-					
6	Phenylcy	clohexyl) - Pyrrolidine, PCPy, PHP;					
7	28.	1-(3-trifluoromethylphenyl) piperazine;					
8	29.	Flunitrazepam;					
9	30.	B-hydroxy-amphetamine;					
10	31.	B-ketoamphetamine;					
11	32.	2,5-dimethoxy-4-nitroamphetamine;					
12	33.	33. 2,5-dimethoxy-4-bromophenethylamine;					
13	34.	34. 2,5-dimethoxy-4-chlorophenethylamine;					
14	35. 2,5-dimethoxy-4-iodoamphetamine;						
15	36.	2,5-dimethoxy-4-iodophenethylamine;					
16	37.	2,5-dimethoxy-4-methylphenethylamine;					
17	38.	2,5-dimethoxy-4-ethylphenethylamine;					
18	39.	2,5-dimethoxy-4-fluorophenethylamine;					
19	40.	2,5-dimethoxy-4-nitrophenethylamine;					
20	41.	2,5-dimethoxy-4-ethylthio-phenethylamine;					
21	42.	2,5-dimethoxy-4-isopropylthio-phenethylamine;					
22	43.	2,5-dimethoxy-4-propylthio-phenethylamine;					
23	44.	2,5-dimethoxy-4-cyclopropylmethylthio-phenethylamine;					
24	45.	2,5-dimethoxy-4-tert-butylthio-phenethylamine;					

2	47.	5-methoxy-N, N-dimethyltryptamine;
3	48.	N-methyltryptamine;
4	49.	A-ethyltryptamine;
5	50.	A-methyltryptamine;
6	51.	N, N-diethyltryptamine;
7	52.	N, N-diisopropyltryptamine;
8	53.	N, N-dipropyltryptamine;
9	54.	5-methoxy-a-methyltryptamine;
10	55.	4-hydroxy-N, N-diethyltryptamine;
11	56.	4-hydroxy-N, N-diisopropyltryptamine;
12	57.	5-methoxy-N, N-diisopropyltryptamine;
13	58.	4-hydroxy-N-isopropyl-N-methyltryptamine;
14	59.	3,4-Methylenedioxymethcathinone (Methylone);
15	60.	3,4-Methylenedioxypyrovalerone (MDPV);
16	61.	3-Methylmethcathinone (Metaphedrone);
17	62.	4-Methylmethcathinone (Mephedrone);
18	63.	4-methoxymethcathinone;
19	64.	4-Fluoromethcathinone;
20	65.	3-Fluoromethcathinone;
21	66.	1-(8-bromobenzo 1,2-b;4,5-b' difuran-4-yl)-2-aminopropane;
22	67.	2,5-Dimethoxy-4-chloroamphetamine;
23	68.	4-Methylethcathinone;
24	69.	Pyrovalerone;

46. 2,5-dimethoxy-4-(2-fluoroethylthio)-phenethylamine;

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 71. 3,4-Methylenedioxy-N-ethylcathinone (Ethylone); 72. B-keto-N-Methylbenzodioxolylbutanamine (Butylone); 73. B-keto-Methylbenzodioxolylpentanamine (Pentylone); 74. Alpha-Pyrrolidinopentiophenone; 75. 4-Fluoroamphetamine; 76. Pentedrone; 77. 4'-Methyl-a-pyrrolidinohexaphenone; 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 80. 1,4-Dibenzylpiperazine; 81. N,N-Dimethylamphetamine; 82. 4-Fluoromethamphetamine; 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine; (25C-NBOMe); 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine; 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine; 86. 1-(4-Fluorophenyl)piperazine; 86. 1-(4-Fluorophenyl)piperazine; 87. Methoxetamine; 88. 3,4-dichloro-N(2-dimethylamino)cyclohexyl]-N- methylbenzamide; 89. N-ethyl hexadrone; 	1	70.	N,N-diallyl-5-methoxytryptamine;			
 Final and the second second	2	71.	71. 3,4-Methylenedioxy-N-ethylcathinone (Ethylone);			
 74. Alpha-Pyrrolidinopentiophenone; 75. 4-Fluoroamphetamine; 76. Pentedrone; 77. 4'-Methyl-a-pyrrolidinohexaphenone; 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 79. 2,5-dimethoxyphenethylamine; 80. 1,4-Dibenzylpiperazine; 81. N,N-Dimethylamphetamine; 82. 4-Fluoromethamphetamine; 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25C-NBOMe); 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25I-NBOMe); 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25B-NBOMe); 86. 1-(4-Fluorophenyl)piperazine; 87. Methoxetamine; 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- methylbenzamide; 	3	72. B-keto-N-Methylbenzodioxolylbutanamine (Butylone);				
 6 75. 4-Fluoroamphetamine; 7 76. Pentedrone; 8 77. 4'-Methyl-a-pyrrolidinohexaphenone; 9 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 10 79. 2,5-dimethoxyphenethylamine; 11 80. 1,4-Dibenzylpiperazine; 12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 	4	73.	B-keto-Methylbenzodioxolylpentanamine (Pentylone);			
 76. Pentedrone; 77. 4'-Methyl-a-pyrrolidinohexaphenone; 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 79. 2,5-dimethoxyphenethylamine; 80. 1,4-Dibenzylpiperazine; 81. N,N-Dimethylamphetamine; 82. 4-Fluoromethamphetamine; 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25C-NBOMe); 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25I-NBOMe); 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine (25B-NBOMe); 86. 1-(4-Fluorophenyl)piperazine; 87. Methoxetamine; 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- methylbenzamide; 	5	74.	Alpha-Pyrrolidinopentiophenone;			
 8 77. 4'-Methyl-a-pyrrolidinohexaphenone; 9 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 10 79. 2,5-dimethoxyphenethylamine; 11 80. 1,4-Dibenzylpiperazine; 12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 	6	75.	4-Fluoroamphetamine;			
 9 78. 2,5-dimethoxy-4-(n)-propylphenethylamine; 10 79. 2,5-dimethoxyphenethylamine; 11 80. 1,4-Dibenzylpiperazine; 12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide; 	7	76.	Pentedrone;			
<pre>10 79. 2,5-dimethoxyphenethylamine; 11 80. 1,4-Dibenzylpiperazine; 12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	8	77.	4'-Methyl-a-pyrrolidinohexaphenone;			
<pre>11 80. 1,4-Dibenzylpiperazine; 12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	9	78.	2,5-dimethoxy-4-(n)-propylphenethylamine;			
<pre>12 81. N,N-Dimethylamphetamine; 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	10	79.	2,5-dimethoxyphenethylamine;			
 13 82. 4-Fluoromethamphetamine; 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide; 	11	80.	1,4-Dibenzylpiperazine;			
 14 83. 4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide; 	12	81.	N,N-Dimethylamphetamine;			
<pre>15 (25C-NBOMe); 16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	13	82.	4-Fluoromethamphetamine;			
<pre>16 84. 4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine 17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	14	83.	4-Chloro-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine			
<pre>17 (25I-NBOMe); 18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	15	(25C-NBO	Me);			
<pre>18 85. 4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine 19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	16	84.	4-Iodo-2,5-dimethoxy-N-(2-methoxybenzyl)phenethylamine			
<pre>19 (25B-NBOMe); 20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	17	(25I-NBO	Me);			
<pre>20 86. 1-(4-Fluorophenyl)piperazine; 21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	18	85.	4-Bromo-2,5-dimethoxy-N-(2-methoxybenzy)phenethylamine			
<pre>21 87. Methoxetamine; 22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;</pre>	19	(25B-NBO	Me);			
22 88. 3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N- 23 methylbenzamide;	20	86.	1-(4-Fluorophenyl)piperazine;			
23 methylbenzamide;	21	87.	Methoxetamine;			
	22	88.	3,4-dichloro-N[2-dimethylamino)cyclohexyl]-N-			
24 89. N-ethyl hexadrone;	23	methylbe	nzamide;			
	24	89.	N-ethyl hexadrone;			

1	90.	Isopropyl-U-47700;
2	91.	Para-fluorobutyrl fentanyl;
3	92.	Para-fluorofentanyl (pFF);
4	93.	Fluoro isobutryrl fentanyl;
5	94.	3-Hydroxy Phencyclidine (PCP);
6	95.	3-methoxy Phencyclidine (PCP);
7	96.	Flualprazolam; or
8	97.	Flubromazolam <u>;</u>
9	98.	Ethyleneoxynitazene;
10	99.	5-Methyl Etodesnitazene;
11	<u>100.</u>	3', 4' Methylenedioxynitazene;
12	101.	N-Pyrrolidino Isotonitazene;
13	<u>102.</u>	N-Desethyl Etonitazene; or
14	<u>103.</u>	Ethylene Etonitazene.
15	D.	Unless specifically excepted or unless listed in a different
16	schedule	, any material, compound, mixture, or preparation which
17	contains	any quantity of the following substances having stimulant
18	or depre	ssant effect on the central nervous system:
19	1.	Fenethylline;
20	2. 1	Mecloqualone;
21	3. 1	N-ethylamphetamine;
22	4. 1	Methaqualone;
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5. Gamma-Hydroxybutyric Acid, also known as GHB, gamma hydroxybutyrate, 4-hydroxybutyrate, 4-hydroxybutanoic acid, sodium
 oxybate, and sodium oxybutyrate;

6. Gamma-Butyrolactone (GBL) as packaged, marketed,
manufactured or promoted for human consumption, with the exception
of legitimate food additive and manufacturing purposes;

7 7. Gamma Hydroxyvalerate (GHV) as packaged, marketed, or
8 manufactured for human consumption, with the exception of legitimate
9 food additive and manufacturing purposes;

10 8. Gamma Valerolactone (GVL) as packaged, marketed, or 11 manufactured for human consumption, with the exception of legitimate 12 food additive and manufacturing purposes;

9. 1,4 Butanediol (1,4 BD or BDO) as packaged, marketed,
manufactured, or promoted for human consumption with the exception
of legitimate manufacturing purposes; or

16 10. N-ethylpentylone.

E. 1. The following industrial uses of Gamma-Butyrolactone,
Gamma Hydroxyvalerate, Gamma Valerolactone, or 1,4 Butanediol are
excluded from all schedules of controlled substances under this
title:

21 a. pesticides,

- 22 b. photochemical etching,
- 23 c. electrolytes of small batteries or capacitors,
- 24 d. viscosity modifiers in polyurethane,

1	e.	surface etching of metal coated plastics,	
2	f.	organic paint disbursements for water soluble inks,	
3	g.	pH regulators in the dyeing of wool and polyamide	
4		fibers,	
5	h.	foundry chemistry as a catalyst during curing,	
6	i.	curing agents in many coating systems based on	
7		urethanes and amides,	
8	j.	additives and flavoring agents in food, confectionary,	
9		and beverage products,	
10	k.	synthetic fiber and clothing production,	
11	1.	tetrahydrofuran production,	
12	m.	gamma butyrolactone production,	
13	n.	polybutylene terephthalate resin production,	
14	ο.	polyester raw materials for polyurethane elastomers	
15		and foams,	
16	p.	coating resin raw material, and	
17	d.	as an intermediate in the manufacture of other	
18		chemicals and pharmaceuticals.	
19	2. At th	e request of any person, the Director of the Oklahoma	
20	State Bureau	of Narcotics and Dangerous Drugs Control may exempt any	
21	other product containing Gamma-Butyrolactone, Gamma Hydroxyvalerate,		
22	Gamma Valerol	actone, or 1,4 Butanediol from being included as a	
23	Schedule I co	ntrolled substance if such product is labeled,	
24			

marketed, manufactured and distributed for legitimate industrial use
 in a manner that reduces or eliminates the likelihood of abuse.

3 3. In making a determination regarding an industrial product,
4 the Director, after notice and hearing, shall consider the
5 following:

6 the history and current pattern of abuse, a. 7 b. the name and labeling of the product, the intended manner of distribution, advertising and 8 с. 9 promotion of the product, and d. other factors as may be relevant to and consistent 10 11 with the public health and safety.

12 4. The hearing shall be held in accordance with the procedures13 of the Administrative Procedures Act.

14 Any material, compound, mixture, or preparation, whether F. 15 produced directly or indirectly from a substance of vegetable origin 16 or independently by means of chemical synthesis, or by a combination 17 of extraction and chemical synthesis, that contains any quantity of 18 the following substances, or that contains any of their salts, 19 isomers, and salts of isomers when the existence of these salts, 20 isomers, and salts of isomers is possible within the specific 21 chemical designation:

22 1. JWH-004;

23 2. JWH-007;

24 3. JWH-009;

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1		4.	JWH-015;
2		5.	JWH-016;
3		6.	JWH-018;
4		7.	JWH-019;
5		8.	JWH-020;
6		9.	JWH-030;
7		10.	JWH-046;
8		11.	JWH-047;
9		12.	JWH-048;
10		13.	JWH-049;
11		14.	JWH-050;
12		15.	JWH-070;
13		16.	JWH-071;
14		17.	JWH-072;
15		18.	JWH-073;
16		19.	JWH-076;
17		20.	JWH-079;
18		21.	JWH-080;
19		22.	JWH-081;
20		23.	JWH-082;
21		24.	JWH-094;
22		25.	JWH-096;
23		26.	JWH-098;
24		27.	JWH-116;
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1	28	3.	JWH-120;
2	29	۶.	JWH-122;
3	3().	JWH-145;
4	31	L.	JWH-146;
5	32	2.	JWH-147;
6	33	3.	JWH-148;
7	34	1.	JWH-149;
8	35	5.	JWH-150;
9	36	5.	JWH-156;
10	37	7.	JWH-167;
11	38	3.	JWH-175;
12	39).	JWH-180;
13	4().	JWH-181;
14	41	L.	JWH-182;
15	42	2.	JWH-184;
16	43	3.	JWH-185;
17	44	1.	JWH-189;
18	45	5.	JWH-192;
19	46	5.	JWH-193;
20	47	7.	JWH-194;
21	48	3.	JWH-195;
22	49).	JWH-196;
23	50).	JWH-197;
24	51	L.	JWH-198;

1	52.	JWH-199;
2	53.	JWH-200;
3	54.	JWH-201;
4	55.	JWH-202;
5	56.	JWH-203;
6	57.	JWH-204;
7	58.	JWH-205;
8	59.	JWH-206;
9	60.	JWH-207;
10	61.	JWH-208;
11	62.	JWH-209;
12	63.	JWH-210;
13	64.	JWH-211;
14	65.	JWH-212;
15	66.	JWH-213;
16	67.	JWH-234;
17	68.	JWH-235;
18	69.	JWH-236;
19	70.	JWH-237;
20	71.	JWH-239;
21	72.	JWH-240;
22	73.	JWH-241;
23	74.	JWH-242;
24	75.	JWH-243;

1	76.	JWH-244;
2	77.	JWH-245;
3	78.	JWH-246;
4	79.	JWH-248;
5	80.	JWH-249;
6	81.	JWH-250;
7	82.	JWH-251;
8	83.	JWH-252;
9	84.	JWH-253;
10	85.	JWH-262;
11	86.	JWH-292;
12	87.	JWH-293;
13	88.	JWH-302;
14	89.	JWH-303;
15	90.	JWH-304;
16	91.	JWH-305;
17	92.	JWH-306;
18	93.	JWH-307;
19	94.	JWH-308;
20	95.	JWH-311;
21	96.	JWH-312;
22	97.	JWH-313;
23	98.	JWH-314;
24	99.	JWH-315;

1	100.	JWH-316;
2	101.	JWH-346;
3	102.	JWH-348;
4	103.	JWH-363;
5	104.	JWH-364;
6	105.	JWH-365;
7	106.	JWH-367;
8	107.	JWH-368;
9	108.	JWH-369;
10	109.	JWH-370;
11	110.	JWH-371;
12	111.	JWH-373;
13	112.	JWH-386;
14	113.	JWH-387;
15	114.	JWH-392;
16	115.	JWH-394;
17	116.	JWH-395;
18	117.	JWH-397;
19	118.	JWH-398;
20	119.	JWH-399;
21	120.	JWH-400;
22	121.	JWH-412;
23	122.	JWH-413;
24	123.	JWH-414;

1	124.	JWH-415;
2	125.	CP-55, 940;
3	126.	CP-47, 497;
4	127.	HU-210;
5	128.	HU-211;
6	129.	WIN-55, 212-2;
7	130.	AM-2201;
8	131.	AM-2233;
9	132.	JWH-018 adamantyl-carboxamide;
10	133.	AKB48;
11	134.	JWH-122 N-(4-pentenyl)analog;
12	135.	MAM2201;
13	136.	URB597;
14	137.	URB602;
15	138.	URB754;
16	139.	UR144;
17	140.	XLR11;
18	141.	A-796,260;
19	142.	STS-135;
20	143.	AB-FUBINACA;
21	144.	AB-PINACA;
22	145.	PB-22;
23	146.	AKB48 N-5-Fluorpentyl;
24	147.	AM1248;

- 1 148. FUB-PB-22;
- 2 149. ADB-FUBINACA;
- 3 150. BB-22;
- 4 151. 5-Fluoro PB-22; or
- 5 152. 5-Fluoro AKB-48.

G. In addition to those substances listed in subsection F of
this section, unless specifically excepted or unless listed in
another schedule, any material, compound, mixture, or preparation
which contains any quantity of a synthetic cannabinoid found to be
in any of the following chemical groups:

11 1. Naphthoylindoles: any compound containing a 3-(1-12 naphthoyl) indole structure with or without substitution at the 13 nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, 14 alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-15 (N-methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-16 2-pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl, 17 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 18 halophenyl group, whether or not further substituted on the indole 19 ring to any extent, and whether or not substituted on the naphthyl 20 ring to any extent. Naphthoylindoles include, but are not limited 21 to: 22 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWHa. 23 200),

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- b. 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201),

1	c. 1-pentyl-3-(1-naphthoyl)indole (JWH-018),
2	d. 1-butyl-3-(1-naphthoyl)indole (JWH-073),
3	e. 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (JWH-081),
4	f. 1-propyl-2-methyl-3-(1-naphthoyl)indole (JWH-015),
5	g. 1-hexyl-3-(1-naphthoyl)indole (JWH-019),
6	h. 1-pentyl-3-(4-methyl-1-naphthoyl)indole (JWH-122),
7	i. 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (JWH-210),
8	j. 1-pentyl-3-(4-chloro-1-naphthoyl)indole (JWH-398),
9	k. 1-pentyl-2-methyl-3-(1-naphthoyl)indole (JWH-007),
10	1. 1-pentyl-3-(7-methoxy-1-naphthoyl)indole (JWH-164),
11	m. 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole
12	(JWH-098),
13	n. 1-pentyl-3-(4-fluoro-1-naphthoyl)indole (JWH-412),
14	o. 1-[1-(N-methyl-2-piperidinyl)methyl]-3-(1-
15	naphthoyl)indole (AM-1220),
16	p. 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole
17	(MAM-2201), or
18	q. 1-(4-cyanobutyl)-3-(1-naphthoyl)indole (AM-2232);
19	2. Naphthylmethylindoles: any compound containing a 1H-indol-
20	3-yl-(1-naphthyl)methane structure with or without substitution at
21	the nitrogen atom of the indole ring by an alkyl, haloalkyl,
22	cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
23	halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
24	morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-

morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or halophenyl group, whether or not further substituted on the indole ring to any extent, and whether or not substituted on the naphthyl ring to any extent. Naphthylmethylindoles include, but are not limited to, (1-pentylindol-3-yl)(1-naphthyl)methane (JWH-175);

6 3. Naphthoylpyrroles: any compound containing a 3-(1-7 naphthoyl)pyrrole structure with or without substitution at the 8 nitrogen atom of the pyrrole ring by an alkyl, haloalkyl, 9 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 10 halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-11 morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, 12 13 phenyl, or halophenyl group, whether or not further substituted on 14 the pyrrole ring to any extent, and whether or not substituted on 15 the naphthyl group to any extent. Naphthoylpyrroles include, but

- 16 are not limited to:
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a. 1-hexyl-2-phenyl-4-(1-naphthoyl)pyrrole (JWH-147),
b. 1-pentyl-5-(2-methylphenyl)-3-(1-naphthoyl)pyrrole (JWH-370),

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c. 1-pentyl-3-(1-naphthoyl)pyrrole (JWH-030), or

d. 1-hexyl-5-phenyl-3-(1-naphthoyl)pyrrole (JWH-147);
4. Naphthylideneindenes: any compound containing a 1-(1naphthylmethylene)indene structure with or without substitution at
the 3-position of the indene ring by an alkyl, haloalkyl,

1 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,

2 halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-

3 morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-4 morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, 5 phenyl, or halophenyl group, whether or not further substituted on 6 the indene group to any extent, and whether or not substituted on 7 the naphthyl group to any extent. Naphthylmethylindenes include, 8 but are not limited to, (1-[(3-pentyl)-1H-inden-1-

9 ylidene)methyl]naphthalene (JWH-176);

10 5. Phenylacetylindoles: any compound containing a 3-11 phenylacetylindole structure with or without substitution at the 12 nitrogen atom of the indole ring by alkyl, haloalkyl, cyanoalkyl, 13 alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-14 (N-methyl-2-piperidinyl) methyl, 2-(4-morpholinyl) ethyl, 1-(N-methyl-15 2-pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl, 16 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 17 halophenyl group, whether or not further substituted on the indole 18 ring to any extent, and whether or not substituted on the phenyl 19 ring to any extent. Phenylacetylindoles include, but are not 20 limited to: 21 1-pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250), a.

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(RCS-8),
c. 1-pentyl-3-(2-chlorophenylacetyl)indole (JWH-203),

1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole

b.

1	d. 1-pentyl-3-(2-methylphenylacetyl)indole (JWH-251),
2	
	e. 1-pentyl-3-(4-methoxyphenylacetyl)indole (JWH-201), or
3	<pre>f. 1-pentyl-3-(3-methoxyphenylacetyl)indole (JWH-302);</pre>
4	6. Cyclohexylphenols: any compound containing a 2-(3-
5	hydroxycyclohexyl)phenol structure with or without substitution at
6	the 5-position of the phenolic ring by an alkyl, haloalkyl,
7	cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl,
8	halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
9	morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
10	morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl,
11	phenyl, or halophenyl group, and whether or not further substituted
12	on the cyclohexyl ring to any extent. Cyclohexylphenols include,
13	but are not limited to:
14	a. 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-
15	hydroxycyclohexyl]-phenol (CP-47,497),
16	b. 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-
17	phenol (cannabicyclohexanol; CP-47,497 C8 homologue),
18	or
19	c. 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-
20	hydroxypropyl)cyclohexyl]-phenol (CP 55, 940);
21	7. Benzoylindoles: any compound containing a 3-(benzoyl)indole
22	structure with or without substitution at the nitrogen atom of the
23	indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
24	cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-

1	2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
2	pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl,
3	(tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
4	halophenyl group, whether or not further substituted on the indole
5	ring to any extent, and whether or not substituted on the phenyl
6	group to any extent. Benzoylindoles include, but are not limited
7	to:
8	a. 1-pentyl-3-(4-methoxybenzoyl)indole (RCS-4),
9	b. 1-[2-(4-morpholinyl)ethyl]-2-methyl-3-(4-
10	methoxybenzoyl)indole (Pravadoline or WIN 48, 098),
11	c. 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM-694),
12	d. 1-pentyl-3-(2-iodobenzoyl)indole (AM-679), or
13	e. 1-[1-(N-methyl-2-piperidinyl)methyl]-3-(2-
14	iodobenzoyl)indole (AM-2233);
15	8. Cyclopropoylindoles: Any compound containing a 3-
16	(cyclopropoyl)indole structure with substitution at the nitrogen
17	atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
18	cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-
19	2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
20	pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl,
21	(tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
22	halophenyl group, whether or not further substituted in the indole
23	ring to any extent and whether or not substituted in the
24	

1 cyclopropoyl ring to any extent. Cyclopropoylindoles include, but 2 are not limited to:

3	a. 1-pentyl-3-(2,2,3,3-tetramethylcyclopropoyl)indole
4	(UR-144),
5	b. 1-(5-chloropentyl)-3-(2,2,3,3-
6	tetramethylcyclopropoyl)indole (5Cl-UR-144), or
7	c. 1-(5-fluoropentyl)-3-(2,2,3,3-
8	<pre>tetramethylcyclopropoyl)indole (XLR11);</pre>
9	9. Indole Amides: Any compound containing a 1H-Indole-3-
10	carboxamide structure with or without substitution at the nitrogen
11	atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
12	cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-
13	2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
14	pyrrolidinyl)methyl, 1-(N-methyl-3- morpholinyl)methyl,
15	(tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
16	halophenyl group, whether or not substituted at the carboxamide
17	group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,
18	cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-
19	1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-
20	dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not
21	further substituted in the indole, adamantyl, naphthyl, phenyl,
22	pyrrole, quninolinyl, or cycloalkyl rings to any extent. Indole
23	Amides include, but are not limited to:

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1	a.	N-(1-adamantyl)-1-pentyl-1H-indole-3-carboxamide
2		(2NE1),
3	b.	N-(1-adamantyl)-1-(5-fluoropentyl-1H-indole-3-
4		carboxamide (STS-135),
5	С.	N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-
6		indole-3-carboxamide (ADBICA),
7	d.	N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-
8		fluoropentyl)-1H-indole-3-carboxamide (5F-ADBICA),
9	e.	N-(naphthalen-1-yl)-1-pentyl-1H-indole-3-carboxamide
10		(NNE1),
11	f.	1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-
12		carboxamide (5F-NNE1),
13	đ.	N-benzyl-1-pentyl-1H-indole-3-carboxamide (SDB-006),
14		or
15	h.	N-benzyl-1-(5-fluoropentyl)-1H-indole-3-carboxamide
16		(5F-SDB-006);
17	10. Indo	le Esters: Any compound containing a 1H-Indole-3-
18	carboxylate s	tructure with or without substitution at the nitrogen
19	atom of the i	ndole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl,
20	cycloalkylmet	hyl, cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-
21	2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
22	pyrrolidinyl)	methyl, 1-(N-methyl-3-morpholinyl)methyl,
23	(tetrahydropy	ran-4-yl)methyl, 1-methylazepanyl, phenyl, or
24	halophenyl gr	oup, whether or not substituted at the carboxylate

1	group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,
2	cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-
3	1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-
4	dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not
5	further substituted in the indole, adamantyl, naphthyl, phenyl,
6	pyrrole, quinolinyl, or cycloalkyl rings to any extent. Indole
7	Esters include, but are not limited to:
8	a. quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate (PB-
9	22),
10	b. quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-
11	carboxylate (5F-PB-22),
12	c. quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-
13	carboxylate (BB-22),
14	d. naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-
15	carboxylate (FDU-PB-22), or
16	e. naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-
17	carboxylate (NM2201);
18	11. Adamantanoylindoles: Any compound containing an
19	adamantanyl-(1H-indol-3-yl)methanone structure with or without
20	substitution at the nitrogen atom of the indole ring by an alkyl,
21	haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl,
22	benzyl, halobenzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-
23	morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
24	morpholinyl)methyl, (tetrahydropyran-4-yl)methyl, 1-methylazepanyl,

phenyl, or halophenyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the adamantyl ring to any extent. Adamantanoylindoles include, but are not limited to:

- a. adamantan-1-yl[1-[(1-methyl-2-piperidinyl)methyl]-1H indol-3-yl]methanone (AM1248), or
- b. adamantan-1-yl-(1-pentyl-1H-indol-3-yl)methanone (AB-001);

9 12. Carbazole Ketone: Any compound containing (9H-carbazole-3yl) methanone structure with or without substitution at the nitrogen 10 11 atom of the carbazole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, halobenzyl, 1-12 13 (N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-14 2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, 15 (tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or 16 halophenyl group, with substitution at the carbon of the methanone 17 group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl, 18 cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-19 1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-20 dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not 21 further substituted at the carbazole, adamantyl, naphthyl, phenyl, 22 pyrrole, quinolinyl, or cycloalkyl rings to any extent. Carbazole 23 Ketones include, but are not limited to, naphthalen-1-yl(9-pentyl-24 9H-carbazol-3-yl)methanone (EG-018);

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1	13. Benzimidazole Ketone: Any compound containing
2	(benzimidazole-2-yl) methanone structure with or without
3	substitution at either nitrogen atom of the benzimidazole ring by an
4	alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
5	cycloalkylethyl, benzyl, halobenzyl, 1-(N-methyl-2-
6	piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
7	pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl,
8	(tetrahydropyran-4-yl)methyl, 1-methylazepanyl, phenyl, or
9	halophenyl group, with substitution at the carbon of the methanone
10	group by an adamantyl, naphthyl, phenyl, benzyl, quinolinyl,
11	cycloalkyl, 1-amino-3-methyl-1-oxobutan-2-yl, 1-amino-3,3-dimethyl-
12	1-oxobutan-2-yl, 1-methoxy-3-methyl-1-oxobutan-2-yl, 1-methoxy-3,3-
13	dimethyl-1-oxobutan-2-yl or pyrrole group, and whether or not
14	further substituted in the benzimidazole, adamantyl, naphthyl,
15	phenyl, pyrrole, quinolinyl, or cycloalkyl rings to any extent.
16	Benzimidazole Ketones include, but are not limited to:
17	a. naphthalen-1-yl(1-pentyl-1H-benzo[d]imidazol-2-
18	l)methanone (JWH-018 benzimidazole analog), or
19	b. (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-
20	yl)(naphthalen-1-yl)methanone (FUBIMINA); and
21	14. Modified by Replacement: any compound defined in this
22	subsection that is modified by replacement of a carbon with nitrogen
23	in the indole, naphthyl, indene, benzimidazole, or carbazole ring.
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1 Η. Any prescription drug approved by the federal Food and Drug 2 Administration under the provisions of Section 505 of the Federal Food, Drug and Cosmetic Act, Title 21 of the United States Code, 3 4 Section 355, that is designated, rescheduled or deleted as a 5 controlled substance under federal law by the United States Drug 6 Enforcement Administration shall be excluded from Schedule I and 7 shall be prescribed, distributed, dispensed or used in accordance 8 with federal law upon the issuance of a notice, final rule or 9 interim final rule by the United States Drug Enforcement 10 Administration designating, rescheduling or deleting as a controlled 11 substance such a drug product under federal law, unless and until 12 the State Board of Pharmacy takes action pursuant to Section 2-201 13 of this title. If the Board of Pharmacy does not take action 14 pursuant to Section 2-201 of this title, the drug product shall be 15 deemed to be designated, rescheduled or deleted as a controlled 16 substance in accordance with federal law and in compliance with the 17 Uniform Controlled Dangerous Substances Act. 18 SECTION 2. This act shall become effective November 1, 2025. 19 20 60-1-10084 GRS 12/31/24

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