1297

TITLE 19 – DEPARTMENT OF HEALTH AND SENIOR SERVICES

Division 30 – Division of Regulation and Licensure Chapter 1 – Controlled Substances

EMERGENCY AMENDMENT

19 CSR 30-1.002 Schedules of Controlled Substances. The department is amending section (1).

PURPOSE: This amendment updates the Schedules of Controlled Substances to be consistent with 21 CFR Part 1308.

EMERGENCY STATEMENT: The United States Department of *Justice Drug Enforcement Administration (DEA) continually* evaluates substances to determine their clinical application and potential for abuse. Based on their evaluation, the DEA issues scheduling actions to place substances in the appropriate controlled substance schedules. The majority of these scheduling actions consist of temporarily and permanently scheduling newlydiscovered illicit substances in Schedule I. Proper scheduling of these substances allow law enforcement to take action to prevent the further distribution of these substances. Scheduling substances in Schedules II-V allows practitioners to be informed about the potential for addiction/abuse of the substances and prescribe the substances appropriately. Section 195.015, RSMo charges the department with similarly controlling substances as they are controlled under federal law. Section 195.015.4 requires the Department of Health and Senior Services to submit emergency rules to the Secretary of State within thirty days of a federal scheduling action to allow for similar inclusion, rescheduling, or deletion of controlled substances with this schedule. While this time frame is difficult to achieve given the various approvals and reviews needed prior to the Department scheduling any rule with the Secretary of State, the Department still acts to effectuate these scheduling actions as quickly as possible. This emergency amendment includes all federal scheduling actions since the last amendment of this rule effective March 20204. This emergency amendment is necessary to protect Missouri's governmental interest in keeping its controlled substances schedules up-to-date as much as practically possible in order to protect its citizens and to aid law enforcement in its prosecution of those who illegally distribute these substances. As a result, the Department of Health and Senior Services finds a compelling governmental interest, which requires this emergency action. The scope of this emergency amendment is limited to the circumstances creating the emergency and complies with the protections extended in the Missouri and United States Constitutions. The Department of Health and Senior Services believes this emergency amendment is fair to all interested persons and parties under the circumstances. Subject to section 536.025, this emergency rule was filed September 24, 2024, becomes effective October 8, 2024, and expires April 5, 2025.

(1) Schedules of Controlled Substances.

(A) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Drug Enforcement Administration (DEA) Controlled Substances Code Number set forth opposite it.

1. Opiates. Unless specifically excepted or unless listed in

another schedule, any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation:

aits is possible within the specific chemic	ai designatio
A. Acetyl-alpha-methylfentanyl	
(N-(1-(1-methyl-2-phenethyl)-	
4-piperidinyl)-N-	
phenylacetamide)	9815
B. Acetylmethadol	9601
C. Acetyl fentanyl (N-(1-	
phenethylpiperidin-4-yl)-	
N-phenylacetamide)	9821
D. N-(1-phenethylpiperidin-	
4-yl)-N-phenylacrylamide,	
its isomers, esters, ethers,	
salts, and salts of isomers,	
esters, and ethers (other	
names: acryl fentanyl,	
acryloylfentanyl)	9811
E. AH-7921(3,4-dichloro-	3011
N-[(1-dimethylamino)	
cyclohexylmethyl]	0551
benzamide)	9551
F. Allylprodine	9602
G. Alphacetylmethadol (except	
levoalphacetylmethadol	
also known as levo-alpha-	
acetylmethadol levothadyl	
acetate or LAAM)	9603
H. Alphameprodine	9604
I. Alphamethadol	9605
J. alpha'-Methyl butyryl fentanyl	
(2-methyl-N-(1-phenethylpiperidin	
-4-yl)-N-phenylbutanamide)	9864
[J.]K. Alpha-methylfentanyl	
[e-jate inplied incomplication of	
(N-1-(alphamethyl-heta-	
(N-1-(alphamethyl-beta-	
phenyl) ethyl-4-piperidyl)	
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl-	
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N-	0014
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine)	9814
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl	9814
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl)	9814
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N-	
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide)	9832
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine	
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide)	9832
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine	9832 9606
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2-	9832 9606
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2-	9832 9606
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)-	9832 9606
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide)	9832 9606 9607
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3-	9832 9606 9607
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name:	9832 9606 9607
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)-	9832 9606 9607
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N-	9832 9606 9607 9830
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide)	9832 9606 9607
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.]Q. N-[1-[2-hydroxy-2-(thiophen-	9832 9606 9607 9830
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.]Q. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]-	9832 9606 9607 9830
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.]L. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.]M. Benzethidine [M.]N. Betacetylmethadol [N.]O. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.]P. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.]Q. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide	9832 9606 9607 9830
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names:	9832 9606 9607 9830
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl)	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.]N. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine [R.JS. Betamethadol	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine [R.JS. Betamethadol [S.JT. beta-Methyl fentanyl	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine [R.JS. Betamethadol [S.JT. beta-Methyl fentanyl (N-phenyl-N-(1-(2-	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine [R.JS. Betamethadol [S.JT. beta-Methyl fentanyl (N-phenyl-N-(1-(2- phenylpropyl)piperidin-4-yl)	9832 9606 9607 9830 9831
phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl- 2-phenylethyl)-4 ((N- propanilido) piperidine) [K.JL. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl) ethyl-4-piperidinyl)-N- phenylpropanamide) [L.JM. Benzethidine [M.JN. Betacetylmethadol [N.JO. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2- phenethyl)-4-piperidinyl)- N-phenylpropanamide) [O.JP. Beta-hydroxy-3- methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)- 3-methyl-4-piperidinyl)-N- phenylpropanamide) [P.JQ. N-[1-[2-hydroxy-2-(thiophen- 2-yl) ethyl]piperidin-4-yl]- N-phenylpropionamide (other names: beta-hydroxythiofentanyl) [Q.JR. Betameprodine [R.JS. Betamethadol [S.JT. beta-Methyl fentanyl (N-phenyl-N-(1-(2-	9832 9606 9607 9830 9831

β-methyl fentanyl)	9856	[QQ.]TT. N-(1-phenethylpiperidin-	
[T.]U. beta'-Phenyl fentanyl		4-yl)-N-phenylfuran-2-	
(N-(1-phenethylpiperidin-4-yl)-		carboxamide (other names:	
N,3-diphenylpropanamide			834
(other names: β'-phenyl		UU. 3-Furanyl fentanyl (N-	
fentanyl; 3-phenylpropanoyl	00.40	(1-phenethylpiperidin-4-yl)-N-	
Fentanyl)	9842		360
[U.]V. Betaprodine	9611		626
[V.]W. Brorphine (1-(1-(4-bromophenyl)			627
ethyl)piperidin-4-yl)-1,3-dihydro-2 <i>H</i> -		[TT.]XX. N-(1-phenethylpiperidin-	
benzo[d]imidazol-2-one)	9098	4-yl)-N-phenylisobutyramide	
[W.]X. Butyryl fentanyl (N-			827
(1-phenethylpiperidin-4-yl)-		[UU.]YY. Isotonitazene (N,N-diethyl-2-(2-	
N-phenylbutyramide)	9822	(4-isopropoxybenzyl)-5-nitro-	
[X.]Y. Clonitazene	9612	1H-benzimidazol-1-yl)	
[Y.]Z. Crotonyl fentanyl ((E)-N-(1-			614
phenethylpiperidin-4-yl)-N-		ZZ. Isovaleryl fentanyl (3-methyl-	
phenylbut-2-enamide)	9844	N-(1-phenethylpiperidin-4-yl)-	
[Z.]AA. N-(1-phenethylpiperidin-		<u> </u>	862
4-yl)-N-			628
Phenylcyclopentanecarboxamide			629
(other name: cyclopentyl			631
fentanyl)	9847	DDD. <i>meta</i> -Fluorofentanyl (<i>N</i> -	
[AA.]BB. Cyclopropyl fentanyl (N-(1-		(3-fluorophenyl)-N-	
phenethylpiperidin-4-yl)-N-		(1-phenethylpiperidin-4-yl)	
phenylcyclopropanecar-			857
boxamide)	9845	EEE. meta-Fluoroisobutyryl fentanyl	
[BB.]CC. Dextromoramide	9613	(N-(3-fluorophenyl)-N-	
[CC.]DD. Diampromide	9615	(1-phenethylpiperidin-4-yl)	
[DD.]EE. Diethylthiambutene	9616		858
[EE.]FF. Difenoxin	9168	[YY.]FFF. Methoxyacetyl fentanyl	
[FF.]GG. Dimenoxadol	9617	(2-methoxy-N-(1-	
[GG.]HH. Dimepheptanol	9618	phenethylpiperidin-4-yl)-	
II. 2',5'-Dimethoxyfentanyl		N-phenylacetamide 98	25
(N-(1-(2,5-dimethoxyphenethyl)		GGG. 2-Methyl AP–237 (1-(2-methyl	
piperidin-4-yl)-N-		-4-(3-phenylprop-2-en-1-yl)	
phenylpropionamide)	9861	piperazin-1-yl)butan-1-one) 960	64
[HH.]JJ. Dimethylthiambutene	9619	[ZZ.]HHH. 4'-Methyl acetyl fentanyl	
[//.]KK. Dioxaphetyl butyrate	9621	(N-(1-(4-methylphenethyl)	
[JJ.]LL. Dipipanone	9622	piperidin-4-yl)-N-	110
[KK.]MM. Ethylmethylthiambutene	9623		319
NN. 2-(2-(4-ethoxybenzyl)-1H-		[AAA.]III. 3-Methylfentanyl (N-(3-	
benzimidazol-1-yl)-N, N-		methyl-1-(2-phenylethyl)-4-	
diethylethan-1-amine (other names:	0=6=	piperidyl)-N-	
etodesnitazene; etazene)	9765	phenylproanamide), its	
[LL.]OO. Etonitazene	9624	optical and geometric	
[MM.]PP. Etoxeridine	9625	isomers, salts, and salts	10
[NN.]QQ. Fentanyl carbamate (ethyl			313
(1-phenethylpiperidin-4-yl)	0054	[BBB.]JJJ. 3-Methylthiofentanyl (N-	
(phenyl)carbamate)	9851	(3-methyl-1-(2-	
[OO.]RR. N-(4-fluorophenyl)-N-		thienyl)ethyl-4-piperidinyl)-	
(1-phenethylpiperidin-4-		N-phenylpropanamide) 98	33
yl)isobutyramide, its		[CCC.]KKK. Metonitazene (N,N-diethyl-	
isomers, esters, ethers,		2-(2-(4-methoxybenzyl)-5-nitro-1H-	
salts, and salts of isomers,		benzimidazol-1-yl)ethan-1-amine) 97	
esters, and ethers (other		[DDD.]LLL. Morpheridine 96.	32
names: 4-fluoroisobutyryl		[EEE.]MMM. MPPP (1-methyl-4-phenyl-4-	
fentanyl, para-			661
fluoroisobutyryl	0004	[FFF.]NNN. MT–45 (1-cyclohexyl-	
fentanyl)	9824	4-(1,2-diphenylethyl)	· ()
[PP.]SS. 2'-Fluoro ortho-		piperazine) (956	,
fluorofentanyl (N-(1-(2-		[GGG.]OOO. Noracymethadol 96	
fluorophenethyl) piperidin-		[HHH.]PPP. Norlevorphanol 96	
4-yl)-N-(2-fluorophenyl)		[///.]QQQ. Normethadone 96	
propionamide (other names:	0055	[JJJ.]RRR. Norpipanone 96	<i>3</i> 6
2'-fluoro 2-fluorofentanyl)	9855	SSS. 2-(4-ethoxybenzyl)-5-nitro-1-	

(2-(pyrrolidin-1-yl)ethyl) -1H- Benzimidazole (other names: N-pyrrolidino etonitazene;		phenethylpiperidin-4- yl)furan-2-carboxamide) <i>[VVV.]</i> FFFF. <i>para</i> -Methoxybutyryl	9854
etonitazepyne)	9758	fentanyl (N-(4-	
[KKK.]TTT. N-(2-fluorophenyl)-2-methoxy-N-(1-	methoxyphenyl)-N-(1-	
phenethylpiperidin-4-		phenethylpiperidin-4-yl)	
yl)acetamide, its isomers,			9837
esters, ethers, salts, and		GGGG. para-Methoxyfuranyl	
salts of isomers, esters,		fentanyl (N-(4-methoxyphenyl)	
and ethers (other name:		-N-(1-phenethylpiperidin-4-yl)	
ocfentanil) `	9838		9859
[LLL.]UUU. ortho-Fluoroacryl fentanyl (N-(2-		HHHH. para-Methylcyclopropyl	
fluorophenyl)-N-(1-		Fentanyl (N-(4-methylphenyl)	
phenethylpiperidin-4-yl)		-N-(1-phenethylpiperidin-4-yl)	
acrylamide)	9852		9865
[MMM.]VVV. ortho-Fluorobutyryl fentanyl		[WWW.]IIII. para-Methylfentanyl	
(N-(2-fluorophenyl)-N-(1-		(N-(4-methylphenyl)-N-	
phenethylpiperidin-4-yl)		(1-phenethylpiperidin-4-yl)	
butyramide (other name:		propionamide (other	
2-fluorobutyryl fentanyl)	9846	name: 4-methylfentanyl)	9817
[NNN.] WWW . ortho-Fluorofentanyl (N-(2-		[XXX.]]]]]. PEPAP (1-(-2-phenethyl)-	
fluorophenyl)-N-(1-		4-phenyl-4-acetoxypiperidine)	9663
phenethylpiperidin-4-yl)		[YYY.]KKKK. Phenadoxone	9637
propionamide); other name:		[ZZZ.]LLLL. Phenampromide	9638
2-fluorofentanyl)	9816	[AAAA.] MMMM . Phenomorphan	9647
XXX. ortho-Fluorofuranyl	3010	[BBBB.]NNNN. Phenoperidine	9641
fentanyl (N-(2-fluorophenyl)		[CCCC.]0000. Phenyl fentanyl (N-(1-	5011
-N-(1-phenethylpiperidin-4-yl)		phenethylpiperidin-4-yl)-	
furan-2-carboxamide)	9863	N-phenylbenzamide	
[OOO.]YYY. ortho-Fluoroisobutyryl	3003	(other name: benzoyl fentanyl)	9841
fentanyl (N-(2-		[DDDD.]PPPP. Piritramide	9642
fluorophenyl)-N-(1-		[EEEE.]QQQQ. Proheptazine	9643
phenethylpiperidin-4-		[FFFF.]RRRR. Properidine	9644
yl)isobutyramide)	9853	[GGGG.]SSSS. Propiram	9649
[PPP.]ZZZ. ortho-Methyl acetylfentanyl	3633	TTTT. N, N-diethyl-2-(5-nitro-2-	JU 1 J
(N-(2-methylphenyl)-N-(1-		(4-propoxybenzyl)-1H-benzimidazol	
phenethylpiperidin-4-yl)		-1-yl)ethan-1-amine (other name:	
acetamide (other name:		Protonitazene)	9759
2-methyl acetylfentanyl)	9848	[HHHH.]UUUU. Racemoramide	9645
[QQQ.]AAAA. ortho-Methyl	3040	[////.]VVVV. N-(1-phenethylpiperidin-4-yl)-	3043
methoxyacetyl fentanyl		<i>N</i> -phenyltetrahydrofuran-	
(2-methoxy-N-(2-		2-carboxamide, its	
methylphenyl)-N-(1-			
		icomore actore athore	
		isomers, esters, ethers,	
phenethylpiperidin-4-yl)		salts, and salts of isomers,	
acetamide (other		salts, and salts of isomers, esters, and ethers (other	
acetamide (other name: 2-methyl		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl	08/13
acetamide (other name: 2-methyl methoxyacetyl	0820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl)	9843
acetamide (other name: 2-methyl methoxyacetyl fentanyl)	9820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.] WWWW . Thiofentany (<i>N</i> -phenyl-	9843
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N-	9820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.] WWWW . Thiofentany (<i>N</i> -phenyl- <i>N</i> -(1-(2-thienyl)ethyl-4-	
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N-(1-phenethylpiperidin-	9820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl-N-(1-(2-thienyl)ethyl-4-piperidinyl)-propanamide	9843 9835
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide	9820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl	
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-	9820	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin-	
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl-N-(1-(2-thienyl)ethyl-4-piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-	
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl)	9820 9826	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other	
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl	9835
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl)	9835 9839
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (other name: parachloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine	9835 9839 9750
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (other name: parachloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-	9826	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine	9835 9839
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide)		salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro-	9835 9839 9750
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N-	9826	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino)	9835 9839 9750
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N-(4-fluorophenyl)-N-(1-(2-	9826	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino) cyclohexyl]-N-	9835 9839 9750 9646
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin- 4-yl)isobutyramide (other name: para- chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4- fluorophenyl)-N-(1- phenethylpiperidin-4- yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N- (4-fluorophenyl)-N-(1-(2- phenethyl)-4-piperidinyl)	9826 9823	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino) cyclohexyl]-N- methylbenzamide)	9835 9839 9750
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N-(4-fluorophenyl)-N-(1-(2-phenethyl)-4-piperidinyl) propanamide	9826	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino) cyclohexyl]-N- methylbenzamide) [OOOO.]CCCCCC. N-(1-phenethylpiperidin-	9835 9839 9750 9646
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin- 4-yl)isobutyramide (other name: para- chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4- fluorophenyl)-N-(1- phenethylpiperidin-4- yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N- (4-fluorophenyl)-N-(1-(2- phenethyl)-4-piperidinyl) propanamide [UUU.]EEEE. para-Fluoro furanyl	9826 9823	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino) cyclohexyl]-N- methylbenzamide) [OOOO.]CCCCC. N-(1-phenethylpiperidin- 4-yl)-N-phenylpentanamide	9835 9839 9750 9646 9547
acetamide (other name: 2-methyl methoxyacetyl fentanyl) [RRR.]BBBB. N-(4-chlorophenyl)-N- (1-phenethylpiperidin-4-yl)isobutyramide (other name: para-chloroisobutyryl fentanyl) [SSS.]CCCC. para-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) [TTT.]DDDD. para-Fluorofentanyl(N-(4-fluorophenyl)-N-(1-(2-phenethyl)-4-piperidinyl) propanamide	9826 9823	salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl) [JJJJ.]WWWW. Thiofentany (N-phenyl- N-(1-(2-thienyl)ethyl-4- piperidinyl)-propanamide [KKKK.]XXXX. Thiofuranyl fentanyl (N-(1-phenethylpiperidin- 4-yl)-N-phenylthiophene- 2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl) [LLLL.]ZZZZ. Tilidine [MMMM.]AAAAA. Trimeperidine [NNNN.]BBBBB. U-47700 (3,4-Dichloro- N-[2-(dimethylamino) cyclohexyl]-N- methylbenzamide) [OOOO.]CCCCCC. N-(1-phenethylpiperidin-	9835 9839 9750 9646

(2-methoxy-2-phenylethyl)		ethanamine (2C-C)	7519
piperazin-1-yl]-1-phenylpropa	n-	M. 2-(4-Ethylthio-2,5-dimethoxyphenyl)	
2-ol)	9873	ethanamine (2C-T-2)	7385
2. Opium derivatives. Unless specifically e		N. 2-(4-Iodo-2,5-dimethoxyphenyl)	
unless listed in another schedule, any of the follow		ethanamine (2C-I)	7518
derivatives, its salts, isomers, and salts of isomer		O. 2-(4-Isopropylthio)-2,5-dimethoxyphenyl)	
the existence of such salts, isomers, and salts o		ethanamine (2C-T-4)	7532
possible within the specific chemical designation:	9319	P. 4-methoxyamphetamine	7411
A. Acetorphine B. Acetyldihydrocodeine	9051	Some trade or other names: 4 amethylphenethylamine; paramethoxyamphetamin	l-methoxy-
C. Benzylmorphine	9052	Q. 5-methoxy-3,4-	ie, rwa,
D. Codeine methylbromide	9070	methylenedioxyamphetamine	7401
E. Codeine-N-Oxide	9053	R. 4-methyl-2,5-dimethoxyamphetamine	7395
F. Cyprenorphine	9054	Some trade and other names: 4-methyl-2, 5- dir	
G. Desomorphine	9055	methylphenethylamine; DOM; and STP;	neurony u
H. Dihydromorphine	9145	S. 3,4-methylenedioxyamphetamine	7400
I. Drotebanol	9335	T. 3,4-methylenedioxymetham-	
J. Etorphine (except hydrochloride salt)	9056	phetamine(MDMA)	7405
K. Heroin	9200	U. 3,4-methylenedioxy-N-	
L. Hydromorphinol	9301	ethylamphetamine (also	
M. Methyldesorphine	9302	known as N-ethylalpha-	
N. Methyldihydromorphine	9304	methyl-3,4 (methylenedioxy)	
O. Morphine methylbromide	9305	phenethylamine, N-ethyl	
P. Morphine methylsulfonate	9306	MDA, MDE, and MDEA)	7404
Q. Morphine-N-Oxide	9307	V. N-hydroxy-3,4-	
R. Myrophine	9308	methylenedioxyamphetamine	
S. Nicocodeine	9309	(also known as N-hydroxy-	
T. Nicomorphine	9312	alpha-methyl-3,4	
U. Normorphine	9313	(methylenedioxy)	
V. Pholcodine	9314	phenethylamine and N-	5 400
W. Thebacon	9315	hydroxy MDA)	7402
3. Hallucinogenic substances. Unless specifica		W. 3,4,5-trimethoxyamphetamine	7390
or unless listed in another schedule, any material, mixture or preparation, which contains any qua		X. 5-MeO-DMT or 5-methoxy-	7431
following hallucinogenic substances or which con		N,N-dimethyltryptamine Y. Alpha-methyltryptamine	7431 7432
its salts, isomers, and salts of isomers whenever the		Z. Bufotenine	7432 7433
of such salts, isomers, and salts of isomers is pos		Some trade and other names: 3-(b-Dimethylan	
the specific chemical designation (For purposes o		5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-ind	
(1)(A)3. of this rule only, the term isomer includes		N-dimethylserotonin; 5-hydroxy-N, N-dimethyltr	
position, and geometric isomers.):	•	mappine;	,,
A. Alpha-ethyltryptamine	7249	AA. Diethyltryptamine	7434
Some trade or other names: etryptamin		Some trade and other names: N, N-Diethyltryptamin	ne; DET;
alpha-ethyl-1H-indole-3-ethenamine; 3-(2-aminob	outyl)indole;	BB. Dimethyltryptamine	7435
alpha-ET; and AET;		Some trade or other names: DMT;	
B. 4-bromo-2,5-dimethoxyamphetamine	7391	CC. 5-methoxy-N,N-diisopropyltryptamine	
Some trade or other names: 4-bromo-2, 5- d	imethoxy-a-	(other name: 5-MeODIPT)	7439
methylphenethylamine; 4-bromo- 2, 5-DMA;	7202	DD. Ibogaine	7260
C. 4-bromo-2,5-dimethoxyphenethylamine D. 2,5-dimethoxyamphetamine	7392 7396	Some trade and other names: 7-Ethyl- 6,6,7,8	
		octahydro-2-methoxy-6, 9-methano-5H-pyrido	[1',2':1,2]
Some trade or other names: 2,5-amethylphenethylamine; 2,5-DMA;	-dimethoxy-	azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide	7315
E. 2,5-dimethoxy-4-ethylamphetamine	7399	FF. Marihuana	7360
Some trade or other names: DOET;	7333	Some trade or other names: marijuana;	7300
F. 2,5-dimethoxy-4-(n)-propylthiophenethyla	amine	GG. Mescaline	7381
(other name: 2C-T-7)	7348	HH. Parahexyl	7374
G. 2-(2,5-Dimethoxy-4-(n)-propylphenyl)	,	Some trade or other names: 3-Hexyl-1- hydro	
ethanamine (2C-P)	7524	tetrahydro-6,6,9-trimethyl- 6H-dibenzo[b,d]pyran; Sy	
H. 2-(2,5-Dimethoxy-4-ethylphenyl)		II. Peyote	7415
ethanamine (2C-E)	7509	Meaning all parts of the plant presently classified by	
I. 2-(2,5-Dimethoxy-4-methylphenyl)		as Lophophora williamsii Lemaire, whether growi	
ethanamine (2C-D)	7508	the seeds thereof; any extract from any part of such	plant; and
J. 2-(2,5-Dimethoxy-4-nitro-		every compound, manufacture, salt, derivative, n	
phenyl) ethanamine (2C-N)	7521	preparation of such plant, its seeds, or extracts;	
K. 2-(2,5-Dimethoxyphenyl)		JJ. N-ethyl-3-piperidyl benzilate	7482
ethanamine (2C-H)	7517	KK. N-methyl-3-piperidyl benzilate	7484
L. 2-(4-Chloro-2,5-dimethoxyphenyl)		LL. Psilocybin	7437
		//	

MM. Psilocyn	7438	EEE. [1-(5-fluoro-pentyl)-	
NN. Tetrahydrocannabinols naturally		1Hindol-3-yl](2,2,3,3-	
plant of the genus Cannabis (cannabis 737		tetramethylcyclopropyl)	
as synthetic equivalents of the substances		methanone (other names:	
cannabis plant or in the resinous extractiv		5-fluoro-UR-144, 5-F-UR-	
and/or synthetic substances, derivatives, and		144, XLR11, 1-(5-fluo-	
both, with similar chemical structure and		ropentyl)-3-(2,2,3,3-	
activity to those substances contained in the	plant, such as the	tetramethylcyclopro-	
following:		poyl)indole)	7011
(I) 1 cis or trans tetrahydrocanna	abinol and their	FFF. N-(1-adamantyl)-1-pentyl-	
optical isomers;		1Hindazole-3-carboxamide	
(II) 6 cis or trans tetrahydrocann	abinol and their	(other names: APINACA, AKB48)	7048
optical isomers;	h:ll :4-	GGG. 2-(4-iodo-2,5-	
(III) 3,4 cis or trans tetrahydrocas optical isomers; and	imadinoi and its	dimethoxyphenyl)-N-(2-	
(IV) Since nomenclature of these	substances is not	methoxybenzyl)ethanamine	
internationally standardized, compounds of		(other names: 251-NBOMe; 2C-I- NBOMe; 25I; Cimbi-5)	7538
regardless of numerical designation of atom		ньоме, 251, Сипог-5) ННН. 2-(4-chloro-2,5-	7556
covered;	inc positions are	dimethoxyphenyl)-N-(2-	
OO. Ethylamine analog of phencyclidi	ne 7455	methoxybenzyl)ethanamine	
Some trade or other names: <i>N</i> -ethyl-1- phenyl		(other names: 25C-NBOMe; 2C-C-NBOMe;	
	henylcyclohexyl)-	25C; Cimbi-82)	7537
ethylamine, cyclohexamine, PCE;	, , , , ,	III. 2-(4-bromo-2,5-	,
PP. Pyrrolidine analog of phencyclidin	e 7458	dimethoxyphenyl)-N-(2-	
Some trade or other names: 1-(1-phenylcyclo	hexyl)-pyrrolidine	methoxybenzyl)ethanamine	
PCPy, PHP;		(other names: 25B-	
QQ. Thiophene analog of phencyclidir		NBOMe; 2C-B-NBOMe;	
Some trade or other names: 1-(1-(2-thier		25B; Cimbi-36)	7536
piperidine, 2-thienyl analog of phencyclidine		JJJ. 4-methyl-N-ethylcathinone	
RR. 1-(1-(2-thienyl)cyclohexyl) pyrrolidi	ne 7473	(other names: 4-MEC; 2-	
Some other names: TCPy;		(ethylamino)-1-(4-	
SS. Salvia divinorum TT. Salvinorin A		methylphenyl)propan-1-one)	1249
UU. 3-Fluoromethcathinone	1233	KKK. 4-methyl-alphapyrrolid-	
VV. 4-Fluoromethcathinone	1238	inopropiophenone,	
WW. Mephedrone, or 4-	1250	(other names: 4-MePPP; MePPP; 4-methyl-	
methylmethcathinone	1248	α -pyrrolidinopropiophenone;	
XX. Methylenedioxy-	12.10	1-(4-methylphenyl)-2-	
pyrovalerone, MDPV, or		(pyrrolidin-1-yl)-propan-1-one)	7498
(1-(1,3-Benzodioxol-5-yl)-		LLL. <i>alpha</i> pyrrolidinopentio-	,
2-(1-pyrrolidinyl)-1-		phenone	
pentanone	7535	(other names: α -PVP; α -	
YY. Methylone, or 3,4-		pyrrolidinovalerophenone;	
Methylenedioxy-		1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one)	7545
methcathinone	7540	MMM. Butylone	
ZZ. Quinolin-8-yl 1-pentyl-		(other names: bk-MBDB; 1-(1,3-	
1Hindole-3-carboxylate	7000	benzodioxol-5-yl)-2-	DE 44
(PB-22; QUPIC) AAA. Quinolin-8-yl 1-(5-	7222	(methylamino)butan-1-one)	7541
fluoropentyl)-1H-indole-		NNN. Pentedrone (other names: α-	
3-carboxylate (5-fluoro-		methylaminovalerophenone;	
PB-22; 5F-PB-22)	7225	2-(methylamino)-1-phenylpentan-1-one)	12/16
BBB. N-(1-amino-3-methyl-1-	7 2 2 3	OOO. Pentylone	12-10
oxobutan-2-yl)-1-		(other names: bk-MBDP; 1-(1,3-	
(4-fluorobenzyl)-1Hindazole-		benzodioxol-5-yl)-2-	
3-carboxamide (AB-FUBINACA)	7012	(methylamino)pentan-1-one)	7542
CCC. N-(1-amino-3, 3-dimethyl-1-		PPP. Naphyrone	
oxobutan-2-yl)-1-pentyl-		(other names:	
1H-indazole-3-carboxamide		naphthylpyrovalerone; 1-	
(ADB-PINACA)	7035	(naphthalen-2-yl)-2-	
DDD. (1-pentyl-1H-indol-3-yl)		(pyrrolidin-1-yl)pentan-1-one)	1258
(2,2,3,3-tetramethylcyclopropyl)		QQQ. <i>alpha</i> -pyrrolidinobutio-	
methanone (other names:		phenone	
UR-144, 1-pentyl-3-(2,2,3,3-		(other names: α-PBP;	7516
tetramethylcyclopro- poyl)indole)	7144	1-phenyl-2-(pyrrolidin-1-yl)butan-1-one) RRR. N-(1-amino-3-methyl-1-	7546
poyijiiidolej	/ 1 '1'1	NNN. 14-(1-ammi0-3-incuryi-1-	

oxobutan-2-yl)-1-		phenylpropan-2-yl)-1 <i>H</i> -	
(cyclohexylmethyl)-		indazole-3-carboxamide	
1H-indazole-3-carboxamide		(other names: 4-CN-	
(other names: AB-CHMINACA)	7031	CUMYLBUTINACA;	
SSS. N-(1-amino-3-methyl-1-		4-cyano-CUMYL-	
oxobutan-2-yl)-1-pentyl-		BUTINACA; 4-CN-	
1Hindazole-3-carboxamide		CUMYLBINACA;	000
(other names:	5000		089
AB-PINACA)	7023	GGGG. methyl 2-(1-(cyclohexylmethyl)-1 <i>H</i> -	
TTT. [1-(5-fluoropentyl)-		indole-3-carboxamido)-3-	
1H-indazol-3-yl](naphthalen-		methylbutanoate	
1-yl)methanone	7024	(other names: MMB-	044
(other names: THJ-2201)	7024	•	044
UUU. N-(1-amino-3,3-dimethyl-		HHHH. 1-(5-fluoropentyl)- <i>N</i> -	
1-oxobutan-2-yl)-1-		(2-phenylpropan-2-yl)-	
(cyclohexylmethyl)-		1H-pyrrolo[2,3-b]	
1H-indazole-3-carboxamide		pyridine-3-carboxamide	085
(other names: MAB-		(Uob
CHMINACA;	7032	IIII. N-ethylpentylone (other	
ADB-CHMINACA)	7032	names: ephylone, 1-(1,3-	
VVV. methyl 2-(1-(5-fluoropentyl)- 1H-indazole-3-carboxamido)-3,3-		benzodioxol-5-yl)-2- (ethylamino)-pentan-1-one)	543
dimethylbutanoate (other names:		[[[]]]. methyl 2-(1-(4-fluorobutyl)-	343
5F–ADB; 5F–MDMB–PINACA)	7034	1H-indazole-3-carboxamido)-3,	
WWW. methyl 2-(1-(5-fluoropentyl)-	7034	3-dimethylbutanoate	
1H-indazole-3-carboxamido)-3-		(4F–MDMB–BINACA,	
methylbutanoate			043
(other names: 5F–AMB)	7033	KKKK. 1-(4-methoxyphenyl)-N-	043
XXX. N-(adamantan-1-yl)-1-(5-	7033	methylpropan-2-amine	
fluoropentyl)-1H-indazole-		(other names: para-	
3-carboxamide		methoxymethamphetamine,	
(other names: 5F–APINACA, 5F–			245
AKB48)	7049	LLLL. ethyl 2-(1-(5-fluoropentyl)-	2-13
YYY. N-(1-amino-3,3-dimethyl-	7043	1 <i>H</i> -indazole-3-carboxamido)-3,3-	
1-oxobutan-2-yl)-1-(4-		dimethylbutanoate	
fluorobenzyl)-1H-indazole-			036
3-carboxamide		MMMM. methyl 2-(1-(5-fluoropentyl)-	000
(other names: ADB–FUBINACA)	7010	1 <i>H</i> -indole-3-carboxamido)-3,3-	
ZZZ. methyl 2-(1-(cyclohexylmethyl)-	, 010	Dimethylbutanoate (other names:	
1H-indole-3-carboxamido)-3,3-			041
dimethylbutanoate		NNNN. N-(adamantan-1-yl)-1-(4-	
(other names: MDMB–CHMICA,		fluorobenzyl)-1 <i>H</i> -indazole-3-	
MMB–CHMINACA)	7042	carboxamide (other names:	
AAAA. methyl 2-(1-(4-fluorobenzyl)-		FUB-AKB48; FUB-APINACA;	
1H-indazole-3-carboxamido)-3,3-			047
dimethylbutanoate		OOOO. 1-(5-fluoropentyl)- <i>N</i> -(2-	
(other names: MDMB–FUBINACA)	7020	phenylpropan-2-yl)-1H-	
BBBB. methyl 2-(1-(4-fluorobenzyl)-1H-		indazole-3-carboxamide (other names:	
indazole-3-carboxamido)-3-			083
methylbutanoate ,		PPPP. (1-(4-fluorobenzyl)-1 <i>H</i> -	
(other names: FUB-AMB, MMB-		indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)	
FUBINACA, AMB-FUBINACA)	(7021)		014
CCCC. 1-(1,3-benzodioxol-	, ,	QQQQ. N-Ethylhexedrone (other names:	
5-yl)-2-(ethylamino)		α-ethylaminohexanophenone; 2-	
propan-1-one (ethylone)	7547	(ethylamino)-1-phenylhexan-1-one) 7	246
DDDD. Naphthalen-1-yl 1-		RRRR. alpha-Pyrrolidinohexanophenone	
(5-fluoropentyl)-1 <i>H</i> -		(other names: α -PHP; α -	
indole-3-carboxylate		pyrrolidinohexanophenone;	
(other names:		1-phenyl-2-(pyrrolidin-1-yl)hexan-1-	
NM2201; CBL2201)	7221	,	544
EEEE. N-(1-amino-3-methyl-		SSSS. 4-Methyl-alpha-ethylaminopentiophenone)
1-oxobutan-2-yl)-1-		(other names: 4-MEAP; 2-(ethylamino)-1-	
(5-fluoropentyl)-1 <i>H</i> -		\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	245
indazole-3-carboxamide		TTTT. 4'-Methyl-alpha-pyrrolidinohexiophenone	
(other name: 5F-AB-PINACA)	7025	(other names: MPHP; 4'-methyl-	
FFFF. 1-(4-cyanobutyl)-N-(2-		alpha-pyrrolidinohexanophenone; 1-	

(4-methylphenyl)-2-(pyrrolidin-1-yl)hexa	n-
1-one)	7446
UUUU. alpha-Pyrrolidinoheptaphenone	
(other names: PV8; 1-phenyl-2-	
(pyrrolidin-1-yl)heptan-1-one)	7548
VVVV. 4'-Chloro-alpha-pyrrolidinovalerophen	one
(other names: 4-chloro-α-PVP; 4'-chloro	
pyrrolidinopentiophenone; 1-(4-	
chlorophenyl)-2-(pyrrolidin-1-yl) pentar	1-
1-one)	7443
WWWW. 2-(ethylamino)-2-	
(3-methoxyphenyl)cyclohexan-	
1-one (methoxetamine, MXE)	7286
XXXX. 1-(1,3-benzodioxol-5-yl)-2-9	
(ethylamino)butan-1-one	
(other names: eutylone;	
bk-EBDB)	7549
YYYY. N-(1-amino-3,3-dimethyl-1-	
oxobutan-2-yl)-1-butyl-1 <i>H</i> -	
indazole-3-carboxamide	
(other name: ADB-BUTINACA)	7027
ZZZZ. 4-methyl-1-phenyl-2-	
(pyrrolidin-1-yl)pentan-1-one	
(other names: α-PiHP;	
alpha-PiHP)	7551
AAAAA. 2-(methylamino)-1-(3-	
methylphenyl)propan-1-one	
(other names: 3–MMC;	
3-methylmethcathinone)	1259
[YYYY.]BBBBB. Synthetic cannabinoids:	Unle

[YYYY.]BBBB. Synthetic cannabinoids: Unless specifically exempted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, or which contains their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(I) Any compound structurally derived from 3-(1-naphthoyl)indole or 1Hindol-3-yl-(1-naphthyl)methane by substitution at the nitrogen atom of the indole ring by 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, whether or not substituted in the naphthyl ring to any extent. Including, but not limited to:

t. Including, but not limited to:	
(a) AM2201, or 1-(5-fluoropentyl)-3-	
(1-naphthoyl)indole	7201
(b) JWH-007, or 1-pentyl-2-methyl-	
3-(1-naphthoyl)indole	
(c) JWH-015, or 1-propyl-2-methyl-3-(1-	
naphthoyl)indole	
(d) JWH-018, or 1-pentyl-	
3-(1-naphthoyl)indole	7118
(e) JWH-019, or 1-hexyl-	
3-(1-naphthoyl)indole	7019
(f) JWH-073, or 1-butyl-	
3-(1-naphthoyl)indole	7173
(g) JWH-081, or 1-pentyl-3-(4-methoxy-1-	
naphthoyl)indole	7081
(h) JWH-098, or 1-pentyl-2-methyl-3-(4-	
methoxy-1-naphthoyl)indole	
(i) JWH-122, or 1-pentyl-3-(4-methyl-1-	
naphthoyl)indole	7122
(j) JWH-164, or 1-pentyl-3-(7-methoxy-1-	
naphthoyl)indole	
(k) JWH-200, or 1-(2-(4-(morpholinyl)ethyl))-
3-(1-naphthoyl)indole	7200

- (l) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole
- (m) JWH-398, or 1-pentyl-3-(4-chloro-1-naphthoyl)indole

7398

6250

(II) Any compound structurally derived from 3-(1-naphthoyl)pyrrole by substitution at the nitrogen atom of the pyrrole ring by 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, whether or not substituted in the naphthyl ring to any extent;

(III) Any compound structurally derived from 1-(1-naphthylmethyl)indene by substitution at the 3-position of the indene ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(*N*-methyl-2- piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indene ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(IV) Any compound structurally derived from 3-phenylacetylindole by substitution at the nitrogen atom of the indole ring with 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, whether or not substituted in the phenyl ring to any extent. Including, but not limited to:

- (a) JWH-201, or 1-pentyl-3-(4-methoxyphenylacetyl)indole
- (b) JWH-203, or 1-pentyl-3-(2-chloropheny-lacetyl)indole 7203(c) JWH-250, or 1-pentyl-
- 3-(2-methoxypheny-lacetyl)indole
- (d) JWH-251, or 1-pentyl-3-(2-methylphenylacetyl)indole
- (e) RCS-8, or 1-(2-cyclohexylethyl)-3-(2methoxypheny-lacetyl)indole 7008
- (V) Any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by 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:

(a) CP 47,497 and homologues, or 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol), where side chain n=5, and homologues where side chain n-4, 6, or 7297.7298

(VI) Any compound containing a 3- (benzoyl)indole structure with substitution at the nitrogen atom of the indole ring by 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:

(a) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole 7694 (b) RCS-4, or 1-pentyl-3-(4-methoxybenzoyl)indole (SR-19 and RCS-4) 7104

(VII) CP 50,556-1, or [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1-yl] acetate;

(VIII) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c] chromen-1-ol:

(IX) HU-211, or Dexanabinol,(6aS,10aS)-9-(hydroxymethyl)- 6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;

(X) Dimethylheptylpyran, or DMHP.

- 4. Depressants. Unless specifically excepted or unless listed in another schedule, any material compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
 - A. Gamma-hydroxybutyric acid and other names GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutonic acid; sodium oxybate;

sodium oxybutryrate2010B. Mecloqualone2572C. Methaqualone2565

- 5. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of isomers:
 - A. Amineptine (7-[(10,11-dihydro-5*H*-dibenzo[*a*,*d*]cyclohepten-5-yl)amino] heptanoic acid)

B. Aminorex 1585

Some trade or other names: aminoxaphen; 2-amino-5-phenyl-

2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine; C. *N*-benzylpiperazine (some

other names: BZP, 1benzylpiperzaine) 7493

- benzylpiperzaine)
 D. Cathinone (Some trade or other names: 2-amino-1-phenyl-1-propanone, alphaaminopropiophenone, 2-aminopropiophenone and norephedrone)
- E. 4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4methyl-5-(4-methylphenyl)-2oxazolamine; 4-methyl-5-(4methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine)

1,3-oxazol-2-amine) 1595
F. Fenethylline 1503
G. Mesocarb (*N*-phenyl-*N* -(3-

(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-yl)carbamimidate)

H. Methcathinone 1237 Some trade or other names: 2-(methylamino)-propiophenone;

Some trade or other names: 2-(methylamino)-propiophenone; alpha-(methylamino) propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-*N*-methylaminopropiophenone; monomethylpropion; ephedrone; *N*-methylcathinone; methylcathinine; AL-464; AL-422; AL-463 and URI 432;

I. Methiopropamine (N-methyl-1-(thiophen-2-yl)propan-2-amine)

1478

1219

1235

1227

J. 4-methoxymethcathinone K. cis-4-methylaminorex

(cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) 1590

L. 4-Methyl-alpha-

pyrrolidinobutiophenone, or MPBP

M. N-ethylamphetamine 1475 N. N,N-dimethylamphetamine 1480

(some other names: *N*,*N*-alpha-trimethylbenzeneethanamine; *N*,*N*-alpha-trimethylphenethylamine)

6. A temporary listing of substances subject to

emergency scheduling under federal law shall include any material, compound, mixture, or preparation which contains any quantity of the following substances:

> A. Fentanyl-related substances, their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers.

9850

9764

9758

2780

- (I) Fentanyl-related substance means any substance not otherwise listed under another Administration Controlled Substance Code Number, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act 21 U.S.C. 355, that is structurally related to fentanyl by one (1) or more of the following modifications:
- (a) Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;
- (b) Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino, or nitro groups;
- (c) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino, or nitro groups:
- (d) Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; and/or
- (e) Replacement of the N-propionyl group by another acyl group.

B. 2-(2-(4-butoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N*, *N*-diethylethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other name: Butonitazene)

(other name: Butonitazene) 9751 [C. 2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-

N, N-diethylethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other names: Etodesnitazene; etazene)]

9765

[D.]C. N, N-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other name: Flunitazene) 9756

(other name: Flunitazene)
[E.JD. N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

(other name: Metodesnitazene)

[F. 2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl) -1H-benzimidazole, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other names: N-pyrrolidino

etonitazene; etonitazepyne)

G. N, N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)1H-benzimidazol-1-yl)ethan-1-amine,
its isomers, esters, ethers, salts, and salts of
isomers, esters and ethers
(other name: Protonitazene)1 9759

(other name: Protonitazene)] [H.]E. 4-(2-chlorophenyl)-2-ethyl-9-

methyl-6*H*-thieno[3,2-*f*][1,2,4]triazolo [4,3-*a*][1,4]diazepine, its salts, isomers, and salts of isomers (Other name: etizolam)

[1.]F. 8-chloro-6-(2-fluorophenyl)-1methyl-4*H*-benzo[*f*][1,2,4]triazolo [4,3-*a*][1,4]diazepine, its salts,

isomers, and salts of isomers (Other name: flualprazolam) [J.]G. 6-(2-chlorophenyl)-1-methyl-8- nitro-4H-benzoz[f][1,2,4]triazolo	2785
[4,3-a][1,4]diazepine, its salts, isomers, and salts of isomers (Other name: clonazolam) [K.]H. 8-bromo-6-(2-fluorophenyl)-1-methyl-4H-benzo[f][1,2,4]triazolo	2786
[4,3-a][1,4]diazepine, its salts, isomers, and salts of isomers (Other name: flubromazolam) [L.]I. 7-chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2 <i>H</i> benzo	2788
[e][1,4]diazepin-2-one, its salts, isomers, and salts of isomers (Other name: diclazepam) J. Methyl 3,3-dimethyl-2-	2789
(1-(pent-4-en-1-yl)-1 <i>H</i> -indazole -3-carboxamido)butanoate, its optical and geometric isomers, salts and salts of isomers	7000
(Other name: MDMB-4en-PINACA) K. Methyl 2-[[1-(4-fluorobutyl) indole-3-carbonyl]amino]-3,3 -dimethyl-butanoate, its optical	7090
and geometric isomers, salts and salts of isomers (Other names: 4F–MDMB–BUTICA; 4F–MDMB–BICA) L. N-(1-Amino-3,3-dimethyl-1-oxobutan -2-yl)-1-(pent-4-en-1-yl)-1H-indazole	7091
-3-carboxamide, its optical and geometric isomers, salts and salts of isomers (Other name: ADB-4en-PINACA) M. 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido	7092
[4,3-b]indol-1-one, its optical and geometric isomers, salts and salts of isomers (Other names: CUMYL–PEGACLO SGT–151) N. Ethyl 2-[[1-(5-fluoropentyl)indole-3	NE; 7093
-carbonyl]amino]-3,3-dimethyl-butanoat its optical and geometric isomers, salts and salts of isomers (Other names:	te,
5F–EDMB–PICA; 5F–EDMB–2201) O. Methyl 2-(1-(4-fluorobenzyl)-1 <i>H</i> -indole-3-carboxamido)-3-methyl butanoate,	7094
its optical and geometric isomers, salts and salts of isomers (Other name: MMB–FUBICA) P. N-ethyl-2-(2-(4-isopropoxybenzyl)-5-	7095
nitro-1H-benzimidazol-1-yl)ethan-1-amino	2,

of isomers, esters and ethers (Other names:
N-piperidinyl etonitazene; etonitazepipne) 9761
7. Khat, to include all parts of the plant presently classified botanically as catha edulis, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seed, or extracts.
7032

its isomers, esters, ethers, salts, and salts

(piperidin-1-yl)ethyl)-1H-benzimidazole,

its isomers, esters, ethers, salts, and salts

N-desethyl isotonitazene)

O. 2-(4-ethoxybenzyl)-5-nitro-1-(2-

of isomers, esters and ethers (Other name:

(B) Schedule II shall consist of the drugs and other substances, by whatever official name, common or usual

- name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Controlled Substances Code Number set forth opposite it.
- 1. Substances, vegetable origin, or chemical synthesis. Unless specifically excepted or unless listed in another schedule, Schedule II shall include any of the following substances whether produced directly or indirectly by extraction from substances of vegetable origin or independently by means of chemical synthesis or by a combination of extraction and chemical synthesis:
- A. Opium and opiate; and any salt, compound, derivative, or preparation of opium or opiate, excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, naloxegol, naloxone, and naltrexone and their respective salts, but including the following:

(I) Raw opium	9	9600
(II) Opium extracts		9610
(III) Opium fluid	!	9620
(IV) Powdered opium		9639
(V) Granulated opium	!	9640
(VI) Tincture of opium	!	9630
(VII) Codeine	•	9050
(VIII) Dihydroetorphine		9334
(IX) Ethylmorphine		9190
(X) Etorphine hydrochloride	!	9059
(XI) Hydrocodone		9193
(XII) Hydromorphone		9150
(XIII) Metopon	!	9260
(XIV) Morphine	!	9300
(XV) Oripavine		9330
(XVI) Oxycodone		9143
(XVII) Oxymorphone		9652
(XVIII) Thebaine		9333

- B. Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in subparagraph (1)(B)1.A. of this rule shall be included in Schedule II, except that these substances shall not include the isoquinoline alkaloids of opium;
 - C. Opium poppy and poppy straw 9650
- D. Coca leaves (9040) and any salt, compound, derivative, or preparation of coca leaves (including cocaine (9041) and ecgonine (9180) and their salts, isomers, derivatives, and salts of isomers and derivatives), and any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of these substances, except that the substances shall not include —
- (I) Decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine; or
 - (II) Ioflupane;
- E. Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid, or powder form which contains the phenanthrene alkaloids of the opium poppy) 9670
- 2. Opiates. Unless specifically excepted or unless in another schedule any of the following opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation, dextrorphan, and levopropoxyphene excepted:

A. Alfentanil	9737
B. Alphaprodine	9010
C. Anileridine	9020
D. Bezitramide	9800
E. Bulk Dextropropoxyphene	
(Non-dosage Forms)	9273

9760

F. Carfentanil	9743
G. Dihydrocodeine	9120
H. Diphenoxylate	9170
I. Fentanyl	9801
J. Isomethadone	9226
K. Levo-alphacetylmethadol	
Some other names: levo-alphaacetylmethadol, levo	methadyl
acetate, LAAM	9648
L. Levomethorphan	9210
M. Levorphanol	9220
N. Metazocine	9240
O. Methadone	9250
P. Methadone-Intermediate,	
4-cyano-2-dimethylamino-	
4,4-diphenyl butane	9254
Q. Moramide-Intermediate, 2-	
methyl-3-morpholino-1,	
1-diphenylpropane-carboxylic acid	9802
R. Oliceridine (N-[(3-methoxythiophen-2-yl)	
methyl] ({2-[(9R)-9-(pyridin-2-yl)-6-oxaspiro	
[4.5]decan-9-yl]ethyl])amine fumarate)	9245
S. Pethidine (Meperidine)	9230
T. Pethidine-Intermediate-A,	
4-cyano-1-methyl-4-phenylpiperidine	9232
U. Pethidine-Intermediate-B,	
ethyl-4-phenylpiperidine-4-carboxylate	9233
V. Pethidine-Intermediate-C, 1-	
methyl-4-phenylpiperidine-	
4-carboxylic acid	9234
W. Phenazocine	9715
X. Piminodine	9730
Y. Racemethorphan	9732
Z. Racemorphan	9733
AA. Remifentanil	9739
BB. Sufentanil	9740
CC. Tapentadol	9780
DD. Thiafentanil	9729
3. Stimulants, Unless specifically excepted or un	less listed

Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system:

A. Amphetamine, its salts, optical isomers,	
and salts of its optical isomers	1100
B. Lisdexamfetamine, its salts, isomers, and	
salts of its isomers	1205
C. Methamphetamine, its salts, isomers, and	
salts of its isomers	1105
D. Phenmetrazine and its salts	1631
E. Methylphenidate	1724

4. Depressants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

	1	
A. Amobarbital		2125
B. Glutethimide		2550
C. Pentobarbital		2270
D. Phencyclidine		7471
E. Secobarbital		2315
5. Hallucinogenic su	bstances:	
A. Nabilone		7379
Another name for nabilo	ne: (±)trans-3-(1,	1- dimethylheptyl)-6,

1-hvdroxv-6.

6a,7,8,10,10a-hexahydro-

dibenzo(b,d) pyran-9-one.

B. Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] in an oral solution in a drug product approved for marketing by the United States Food and Drug Administration. (7365)

6. Immediate precursors. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

A. Immediate precursor to amphetamine methamphetamine:

(I) Phenylacetone

Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;

B. Immediate precursors to phencyclidine (PCP):

(I) 1-phenylcyclohexylamine 7460 (II) 1-piperidinocyclohexanecarbonitrile (PCC) 8603 C. Immediate precursor to fentanyl:

(I) 4-anilino-N-phenethyl-4-

piperidine (ANPP) 8333 (II) N-phenyl-N-(piperidin-

4-yl)propionamide (norfentanyl)

8366 7. Any material, compound, mixture, or preparation which contains any quantity of the following alkyl nitrites:

A. Amyl nitrite;

B. Butyl nitrite.

(C) Schedule III shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.

1. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers (whether optical, position, or geometric), and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

A. Those compounds, mixtures, or preparations in dosage unit form containing any stimulant substances listed in Schedule II which compounds, mixtures, or preparations were listed on August 25, 1971, as excepted compounds under 21 CFR 308.32 and any other drug of the quantitive composition shown in that list for those drugs or which is the same except that it contains a lesser quantity of controlled substances 1405

B. Benzphetamine	1228
C. Chlorphentermine	1645
D. Clortermine	1647
E. Phendimetrazine	1615

2. Depressants. Unless specifically excepted or unless listed in another schedule, any material compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system:

> A. Any compound, mixture, or preparation containing – (I) Amobarbital 2126 2316

(II) Secobarbital (III) Pentobarbital 2271

or any salt thereof and one (1) or more other active medicinal ingredients which are not listed in any schedule;

B. Any suppository dosage form containing – (I) Amobarbital

2126 (II) Secobarbital 2316 (III) Pentobarbital

or any salt of any of these drugs and approved by

6-dimethyl-9H-

the Food and Drug Administration for marketing only as a suppository;

- C. Any substance which contains any quantity of a derivative of barbituric acid or any salt thereof
 D. Chlorhexadol
 2510
 - E. Embutramide 2020
- F. Any drug product containing gamma hydroxybutyric acid, including its salts, isomers, and salts of isomer, for which an application is approved under section 505 of the Federal Food, Drug, and Cosmetic Act;
 - G. Ketamine, its salts, isomer, and salts of isomers (some other names for ketamine: (±)-2-(2-chlorophenyl)-2-(methylamino)-

cyclohexanone) 7285
H. Lysergic acid 7300
I. Lysergic acid amide 7310
J. Methyprylon 2575
K. Poramonal, and its salts isomers, and salts

K. Perampanel, and its salts, isomers, and salts

of isomers 2261
L. Sulfondiethylmethane 2600
M. Sulfonethylmethane 2605
N. Sulfonmethane 2610

O. Tiletamine and zolazepam

or any salt thereof 7295

Some trade or other names for a tiletaminezolazepam combination product: Telazol.

Some trade or other names for tiletamine: 2- (ethylamino)-2-(2-thienyl)-cyclohexanone.

Some trade or other names for zolazepam: 4-(2-fluorophenyl)-6-8-dihydro-1,3,8- trimethylpyrazolo-(3,4-e) (1,4)-diazepin-7(1H)-one, flupyrazapon.

3. Nalorphine 940

4. Narcotics drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof:

A. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium 9803

- B. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts
- C. Not more than one and eight tenths grams (1.8gm) of dihydrocodeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9807
- D. Not more than three hundred milligrams (300 mg) of ethylmorphine per one hundred milliliters (100 mL) or not more than fifteen milligrams (15 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9808
- E. Not more than five hundred milligrams (500 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm) or not more than twenty-five milligrams (25 mg) per dosage unit, with one (1) or more active nonnarcotic ingredients in recognized therapeutic amounts 9809
- F. Not more than fifty milligrams (50 mg) of morphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm), with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9810
 - 5. Any material, compound, mixture, or preparation

containing any of the following narcotic drugs or their salts, as set forth below:

A. Buprenorphine 906

- 6. Anabolic steroids. Unless specifically excepted or unless listed in another schedule, any substance meeting the definition of anabolic steroid as set forth in 21 CFR 1300.01, of the August 1, 2023, *Code of Federal Regulations*, as published by the Office of the Federal Register, National Archives and Records Administration, 700 Pennsylvania Avenue, Washington, DC 20408, which is hereby incorporated by reference and does not include later amendments or additions, including any material, compound, mixture or preparation containing any quantity of the following substances, including its salts, esters, and ethers (4000):
 - A. 5α -androstan-3,17-dione;
 - B. 5α -androstan-3,6,17-trione;
 - C. 1-androstenediol (3β , 17β -dihydroxy- 5α -androst-1-ene);
 - D. 1-androstenediol (3α ,17 β -dihydroxy- 5α -androst-1-ene);
 - E. 4-androstenediol (3β,17β-dihydroxy-androst-4-ene);
 - F. 5-androstenediol (3β ,17 β -dihydroxy-androst-5-ene);
 - G. 1-androstenedione (5α -androst-1-en-3,17-dione);
 - H. 4-androstenedione (androst-4-en-3,17-dione);
 - I. 5-androstenedione (androst-5-en-3,17-dione);
- J. bolasterone (7 α ,17 α -dimethyl-17 β -hydroxyandrost-4-en-3-one);
 - K. boldenone (17β-hydroxyandrost-1,4-diene-3-one);
 - L. boldione (androsta-1,4-diene-3,17-dione);
 - M. 6-bromo-androsta-1.4-diene-3.17-dione:
 - N. 6-bromo-androstan-3,17-dione;
- O. calusterone (7 β ,17 α -dimethyl-17 β -hydroxyandrost-4-en-3-one);
 - P. 4-chloro- 17α -methyl-androsta-1,4-diene-3,17 β -diol;
 - Q. 4-chloro-17α-methyl-androst-4-ene-3β,17β-diol;
 - R. 4-chloro- 17α -methyl- 17β -hydroxy-androst-4-en-3-one;
- S. 4-chloro-17 α -methyl-17 β -hydroxy-androst-4-ene-3,11-dione;
 - T. clostebol (4-chloro-17β-hydroxy-androst-4-en-3-one);
- U. dehydrochloromethyltestosterone (4-chloro-17 β -hydroxy-17 α -methyl-androst-1,4-dien-3-one);
- V. desoxymethyltestosterone (17α -methyl- 5α -androst-2-en- 17β -ol) (a.k.a. "madol");
- W. 4-dihydrotestosterone (17 β -hydroxy-androstan-3-one);
- X. Δ 1-dihydrotestosterone (a.k.a. "1-testosterone") (17 -hydroxy- 5α -androst-1-en-3-one);
 - Y. 3β , 17β -dihydroxy- 5α -androstane;
 - Z. 3α ,17 β -dihydroxy- 5α -androstane;
 - AA. 2α , 17α -dimethyl- 17β -hydroxy- 5β -androstan-3-one;
- BB. drostanolone (17 β -hydroxy-2 α -methyl-5 α -androstan-3-one);
 - CC. 2α , 3α -epithio- 17α -methyl- 5α -androstan- 17β -ol;
 - DD. estra-4,9,11-triene-3,17-dione;
 - EE. 13β -ethyl- 17β -hydroxygon-4-en-3-one;
 - FF. ethylestrenol (17α-ethyl-17β-hydroxyestr-4-ene);
- GG. fluoxymesterone (9-fluoro- 17α -methyl- 11β ,17 β -dihydroxyandrost-4-en-3-one);
- HH. formebolone (2-formyl-17 α -methyl-11 α ,17 β -dihydroxyandrost-1,4-dien-3-one);
- II. furazabol (17 α -methyl-17 β -hydroxyandrostano[2,3-c] furazan):
 - []. [3,2-c]furazan-5α-androstan-17β-ol;
 - KK. 18a-homo-3-hydroxy-estra-2,5(10)-dien-17-one;
- LL. 4-hydroxy-19-nortestosterone (4,17 β -dihydroxy-estr-4-en-3-one);
 - MM. 4-hydroxy-androst-4-ene-3,17-dione;
 - NN. 17β-hydroxy-androstano[2,3-d]isoxazole;

OO. 17β-hydroxy-androstano[3,2-c]isoxazole;

PP. 3β-hydroxy-estra-4,9,11-trien-17-one;

QQ. 4-hydroxytestosterone (4,17 β -dihydroxy-androst-4-en-3-one);

RR. mestanolone (17 α -methyl-17 β -hydroxy-5 α -androstan-3-one);

SS. mesterolone (1α -methyl- 17β -hydroxy- 5α -androstan-3-one);

TT. methandienone (17 α -methyl-17 β -hydroxyandrost-1,4-dien-3-one);

UU. methandriol (17 α -methyl-3 β ,17 β -dihydroxyandrost-5-ene);

VV. methasterone (2 α ,17 α -dimethyl-5 α -androstan-17 β -ol-3-one or 2 α ,17 α -dimethyl-17 β -hydroxy-5 α -androstan-3-one);

WW. methenolone (1-methyl-17 β -hydroxy-5 α -androst-1-en-3-one);

XX. 17α -methyl-androsta-1,4-diene-3,17 β -diol;

YY. 17α-methyl-5α-androstan-17β-ol;

ZZ. 17α-methyl-androstan-3-hydroxyimine-17 -ol;

AAA. 6α -methyl-androst-4-ene-3,17-dione;

BBB. 17α-methyl-androst-2-ene-3,17β-diol;

CCC. 17α -methyl- 3β , 17β -dihydroxy- 5α -androstane;

DDD. 17α -methyl- 3α , 17β -dihydroxy- 5α -androstane;

EEE. 17α-methyl-3β,17β-dihydroxyandrost-4-ene;

FFF. 17α -methyl-4-hydroxynandrolone (17 α -methyl-4-hydroxy-17 β -hydroxyestr-4-en-3-one);

GGG. methyldienolone (17 α -methyl-17 β -hydroxyestra-4.9(10)-dien-3-one);

HHH. 17α -methyl- Δ 1-dihydrotestosterone (17β -hydroxy- 17α -methyl- 5α -androst-1-en-3-one) (a.k.a. "17- α -methyl-1-testosterone");

III. me'thyltestosterone (17 α -methyl-17 β -hydroxyandrost-4-en-3-one);

JJJ. methyltrienolone (17α-methyl-17β-hydroxyestra-4,9,11-trien-3-one);

KKK. mibolerone (7α , 17α -dimethyl- 17β -hydroxyestr-4-en-3-one);

LLL. nandrolone (17β-hydroxyestr-4-en-3-one);

MMM. 19-nor-4-androstenediol (3 β ,17 β -dihydroxyestr-4-ene);

NNN. 19-nor-4-androstenediol $(3\alpha,17\beta$ -dihydroxyestr-4-ene);

OOO. 19-nor-5-androstenediol (3 β ,17 β -dihydroxyestr-5-ene);

PPP. 19-nor-5-androstenediol (3 α ,17 β -dihydroxyestr-5-ene);

QQQ. 19-nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione);

RRR. 19-nor-4-androstenedione (estr-4-en-3,17-dione);

SSS. 19-nor-5-androstenedione (estr-5-en-3,17-dione);

TTT. norbolethone (13β , 17α -diethyl- 17β -hydroxygon-4-en-3-one);

UUU. norclostebol (4-chloro-17 β -hydroxyestr-4-en-3-one);

VVV. norethandrolone (17 α -ethyl-17 β -hydroxyestr-4-en-3-one);

WWW. normethandrolone (17 α -methyl-17 β -hydroxyestr-4-en-3-one);

XXX. oxandrolone (17 α -methyl-17 β -hydroxy-2-oxa-5 -androstan-3-one);

YYY. oxymesterone (17 α -methyl-4,17 β -dihydroxyandrost-4-en-3-one);

ZZZ. oxymetholone (17 α -methyl-2-hydroxymethylene-17 -hydroxy-5 -androstan-3-one);

AAAA. prostanozol (17 β -hydroxy-5 α -androstano[3,2-c] pyrazole or [3,2-c]pyrazole-5 α -androstan-17 β -ol);

BBBB. [3,2-c]pyrazole-androst-4-en-17β-ol;

CCCC. stanozolol (17 α -methyl-17 β -hydroxy-5 α -androst-2-eno[3,2-c]-pyrazole);

DDDD. stenbolone (17 β -hydroxy-2-methyl-5 α -androst-1-en-3-one);

EEEE. testolactone (13-hydroxy-3-oxo-13,17-secoandrosta-1,4-dien-17-oic acid lactone);

FFFF. testosterone (17β-hydroxyandrost-4-en-3-one);

GGGG. tetrahydrogestrinone (13 β ,17 α -diethyl-17 β -hydroxygon-4,9,11-trien-3-one); and

HHHH. trenbolone (17β-hydroxyestr-4,9,11-trien-3-one).

7. Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a United States Food and Drug Administration approved drug product 7369

(Some other names for dronabinol: (6 Rtrans)- $6a,7,8,10\alpha$ -tetrahydro-6.6.9-trimethyl-3-pentyl-6H-dibenzo (b,d) pyran-1-ol, or (-) -delta-9-(trans)-tetrahydrocannabinol.)

(D) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this subsection. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.

1. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof:

A. Not more than one milligram (1 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit 9167

B. Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane) 9278

C. 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol, its salts, optical and geometric isomers, and salts of these isomers (including tramadol) 9752

D. Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following limited quantities of narcotic drugs or salts thereof, which shall include one (1) or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:

(I) Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

(II) Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm); or

(III) Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm).

2. Depressants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

1	J
A. Alfaxalone	2731
B. Alprazolam	2882
C. Barbital	2145
D. Brexanolone	2400
E. Bromazepam	2748
F. Camazepam	2749
G. Carisoprodol	8192
H. Chloral betaine	2460
I. Chloral hydrate	2465

	J. Chlordiazepoxide	2744	
	K. Clobazam	2751	
	L. Clonazepam	2737	
	M. Clorazepate	2768	
	N. Clotiazepam	2752	
	O. Cloxazolam	2753	
	P. Daridorexant	2410	
	Q. Delorazepam	2754	
	R. Diazepam	2765	
	S. Dichloralphenazone	2467	
	T. Estazolam	2756	
	U. Ethchlorvynol	2540	
	V. Ethinamate	2545	
	W. Ethyl loflazepate	2758	
	X. Fludiazepam	2759	
	Y. Flunitrazepam	2763	
	Z. Flurazepam	2767	
	AA. Fospropofol	2138	
	BB. Halazepam	2762	
	CC. Haloxazolam	2771	
	DD. Ketazolam	2772	
	EE. Lemborexant	2245	
	FF. Loprazolam	2773	
	GG. Lorazepam	2885	
	HH. Lormetazepam	2774	
	II. Mebutamate	2800	
	JJ. Medazepam	2836	
	KK. Meprobamate	2820	
	LL. Methohexital	2264	
	MM. Methylphenobarbital (Mephobarbital)	2250	
	NN. Midazolam	2884	
	OO. Nimetazepam	2837	
	PP. Nitrazepam	2834	
	QQ. Nordiazepam	2838	
	RR. Oxazepam	2835	
	SS. Oxazolam	2839	
	TT. Paraldehyde	2585	
	UU. Petrichloral	2591	
	VV. Phenobarbital	2285	
	WW. Pinazepam	2883	
	XX. Prazepam	2764	
	YY. Quazepam	2881	
	ZZ. Remimazolam	2846	
	AAA. Suvorexant	2223	
	BBB. Temazepam	2925	
	CCC. Tetrazepam	2886	
	DDD. Triazolam	2887	
	EEE. Zaleplon	2781	
	FFF. Zolpidem	2783	
	GGG. Zopiclone	2784	
	HHH. Zuranolone	2420	
3.	Lorcaserin. Any material, compound,	mixture,	

3. Lorcaserin. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers is possible:

A. Lorcaserin 1625

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

A. Cathine ((+)-norpseudoephedrine)	1230
B. Diethylpropion	1610
C. Fencamfamin	1760
D. Fenproporex	1575

E. Mazindol	1605
F. Mefenorex	1580
G. Modafinil	1680
H. Pemoline (including organometallic	
complexes and chelates thereof)	1530
I. Phentermine	1640
J. Pipradrol	1750
K. Serdexmethylphenidate	1729
L. Sibutramine	1675
M. Solriamfetol (2-amino-3-phenylpropyl	
carbamate; benzenepropanol, beta-	
amino-, carbamate (ester))	1650

N. SPA (-)-1-dimethylamino-1,2-diphenylethane 1635

5. Other substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts:

A. Pentazocine 9709

B. Butorphanol (including its optical isomers) 9720

C. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropyl] [(1S)-1-(4-phenyl-1 H-imidazol-2-yl) ethyl]amino]methyl]-2-methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and salts of isomers

6. Ephedrine. Any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system including their salts, isomers, and salts of isomers:

A. Ephedrine or its salts, optical isomers, or salts of optical isomers as the only active medicinal ingredient or contains ephedrine or its salts, optical isomers, or salts of optical isomers and therapeutically insignificant quantities of another active medicinal ingredient.

(E) Schedule V shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this subsection.

- 1. Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as follows, which shall include one (1) or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:
- [A. Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);]
- [B.]A. Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);
- [C.]B. Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);
- [D.]C. Not more than two and five-tenths milligrams (2.5 mg) of diphenoxylate and not less than twenty-five micrograms (25 mg) of atropine sulfate per dosage unit;
- [E.]D. Not more than one hundred milligrams (100 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm); and
- [F.JE. Not more than five-tenths milligram (0.5 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit.
- 2. Stimulants. Unless specifically exempted or excluded or unless listed in another schedule, any material, compound,

mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system including its salts, isomers, and salts of isomers:

A. Pyrovalerone 1485

- 3. Any compound, mixture, or preparation containing any detectable quantity of pseudoephedrine or its salts or optical isomers, or salts of optical isomers or any compound, mixture, or preparation containing any detectable quantity of ephedrine or its salts or optical isomers, or salts of optical isomers if the drug preparations are starch-based solid dose forms, if such preparations are sold over the counter without a prescription. The following drug preparations containing ephedrine and pseudoephedrine are not scheduled controlled substances:
 - A. Drug preparations in liquid form; and
- B. Drug preparations that require a prescription in order to be dispensed.
- 4. Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts:

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A. Ezogabine [N-[2-amino-4(4-	
fluorobenzylamino)-phenyl]-	
carbamic acid ethyl ester]	2779
B. Ganaxolone (3α-hydroxy-3β-methyl-	
5 -pregnan-20-one)	2401
C. Lacosamide [(R)-2-acetoamido-N-benzyl-	
3-methoxy-propionamide]	2746
D. Pregabalin [(S)-3-(aminomethyl)-5-	
methylhexanoic acid]	2782
E. Brivaracetam ((25)-2-[(4R)-	
2-oxo-4-propylpyrrolidin-1-	
yl]butanamide) (also referred	
to as BRV; UCB-34714; Briviact)	2710
F. Lasmiditan [2,4,6-trifluoro- <i>N</i> -(6-(1-	
methylpiperidine-4-carbonyl) pyridine-2-	
yl-benzamide]	2790
G. Cenobamate ([(1R)-1-(2-	
chlorophenyl)-2-(tetrazol-	
2-yl)ethyl] carbamate; 2H-	
tetrazole-2-ethanol, alpha-(2-	
chlorophenyl)-, carbamate	
(ester), (alphaR)-; carbamic	
acid (R)-(+)-1-(2-chlorophenyl)-2-(2H-tetrazol-	
2-yl)ethyl ester)	2720

AUTHORITY: section 195.015, RSMo Supp. [2023] 2024, and section 195.195, RSMo 2016. Material found in this rule previously filed as 19 CSR 30-1.010. Original rule filed April 14, 2000, effective Nov. 30, 2000. For intervening history, please consult the Code of State Regulations. Emergency amendment filed Sept. 24. 2024, effective Oct. 8, 2024, expires April 5, 2025. An emergency amendment and proposed amendment covering the same material will be published in the Nov. 1, 2024, issue of the Missouri Register.

PUBLIC COST: This emergency amendment will not cost state agencies or political subdivisions more than five hundred dollars (\$500) in the time the emergency is effective.

PRIVATE COST: This emergency amendment will not cost private entities more than five hundred dollars (\$500) in the time the emergency is effective.