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 newly-discovered 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. This section also now (as of August 28, 2020) states that the Department shall 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 in 2019. The scope of this emergency rule 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 rule is fair to all interested persons and parties under the circumstances. Subject to section 536.025, this emergency rule was filed Oct. 30, 2020, becomes effective Nov. 16, 2020, and expires May 14, 2021.

(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:
         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-N-[1-(1-dimethylamino)cyclohexylmethyl]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-methylfentanyl (N-1-(alpha-methyl-betaphenyl)ethyl-4-piperidyl)propionanilide; 1-(1-methyl-2-phenylethyl)-4-((N-propanilido) piperidine) 9814
K. Alpha-methylthiofentanyl (N-(1-methyl-2-(2-thienyl)ethyl-4-piperidinyl)-N-phenylpropanamide) 9832
L. Benzethidine 9606
M. Betacetylmethadol 9607
N. Beta-hydroxyfentanyl (N-(1-(2-hydroxy-2-phenethyl)-4-piperidinyl)-N-phenylpropanamide) 9830
O. Beta-hydroxy-3-methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl)-N-phenylpropanamide) 9831
P. N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide
   (Other names: beta-hydroxythiofentanyl)  9836

[P] Q. Betameprodine  9608
[Q] R. Betamethadol  9609
[R] S. Betaprodine  9611
[S] T. Clonitazene  9612

U. Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide)  9845

[T] V. Dextromoramide  9613
[U] W. Diampromide  9615
[V] X. Diethylthiambutene  9616
[W] Y. Difenoxin  9168
[X] Z. Dimenoxadol  9617
[Y] AA. Dimepheptanol  9618
[Z] BB. Dimethylthiambutene  9619

[AA] CC. Dioxaphetyl butyrate  9621
[BB] DD. Dipipanone  9622
[CC] EE. Ethylmethylthiambutene  9623
[DD] FF. Etonitazene  9624

[EE] GG. Etoxeridine  9625

[FF] HH. N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other names: 4-fluoroisobutyryl fentanyl, parafluoroisobutyryl fentanyl)  9824

[GG] II. N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (Other names: furanyl fentanyl)  9834

[HH] JJ. Furethidine  9626
[JJ] KK. Hydroxypethidine  9627
[KK] LL. Ketobemidone  9628
[LL] MM. Levomoramide  9629
[MM] NN. Levophenacylmorphan  9631

OO. Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide)  9825

[MM] PP. 3-Methylfentanyl (N-(3-
methyl-1-(2-phenylethyl)-4-piperidyl)-N-phenylpropanamide), its optical and geometric isomers, salts, and salts of isomers 9813

[NN/QQ]. 3-Methylthiofentanyl (N-((3-methyl-1-(2-thienyl)ethyl-4-piperidinyl)-N-phenylpropanamide) 9833

[OO/RR]. Morpheridine 9632

[PP/SS]. MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) 9661

[QQ/TT]. MT–45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine) (9560)

[RR/UU]. Noracymethadol 9633

[SS/VV]. Norlevorphanol 9634

[TT/WW]. Normethadone 9635

[UU/XX]. Norpipanone 9636

[VV/YY]. N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: ofcfentanil) 9838

ZZ. ortho-Fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide); other name: 2-fluorofentanyl) 9816

AAA. para-Fluorobutyrylfentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) 9823

[WW/BBB]. Para-fluorofentanyl(N-(4-fluorophenyl)-N-(1-(2-phenethyl)-4-piperidinyl) propanamide 9812

[XX/CCC]. PEPAP (1-(2-phenethyl)-4-phenyl-4-acetoxyxypiperidine) 9663

[YY/DDD]. Phenadoxone 9637

[ZZ/EEE]. Phenampromide 9638

[AAA/FFF]. Phenomorphan 9647
[BBB]GGG. Phenoperidine 9641
[CCC/HHH. Piriramide 9642
[DDD/III. Proheptazine 9643
[EEE]JJJ. Properidine 9644
[FFF]KKK. Propiram 9649
[GGG]LLL. Racemoramide 9645
[HHH/MMM. N-(1-phenethylpiperidin-4-yl)-N-
phenyltetrahydrofuran-2-carboxamide, its
isomers, esters, ethers,
salts and salts of isomers,
esters and ethers (Other
name: tetrahydrofuranyl
fentany) 9843
[III]NNN. Thiofentany (N-phenyl-N-
(1-(2-thienyl)ethyl-4-
piperidinyl)-propanamide 9835
[JJJ]OOO. Tilidine 9750
KKK/PPP. Trimeperidine 9646

2. Opium derivatives. Unless specifically
excepted or unless listed in another schedule,
any of the following opium derivatives,
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. Acetorphine 9319
B. Acetyldihydrocodeine 9051
C. Benzyldimorphine 9052
D. Codeine methylbromide 9070
E. Codeine-N-Oxide 9053
F. Cyprenorphine 9054
G. Desomorphine 9055
H. Dihydromorphine 9145
I. Drotebanol 9335
J. Etorphine (except
hydrochloride salt) 9056
K. Heroin 9200
L. Hydromorphinol 9301
M. Methyldeesorphine 9302
N. Methylene codeine methylbromide 9304
O. Morphine methylbromide 9305
P. Morphine methylsulfonate 9306
Q. Morphine-N-Oxide 9307
R. Myrophine 9308
S. Nicocodeine
T. Nicomorphine
U. Normorphine
V. Pholcodine
W. Thebacon

3. Opiate Similar Synthetic Substances.
Substances scheduled by the United States Drug Enforcement Administration as substances that share a pharmacological profile similar to fentanyl, morphine, and other synthetic opioids, unless specifically excepted or unless listed in another schedule. These substances are:

A. Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide) 9822
B. U–47700 (3,4-Dichloro-N-[2-(dimethylamino)cyclohexyl]-Nmethylbenzamide) 9547

4. Hallucinogenic substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation, which contains any quantity of the following hallucinogenic substances or which contains any of 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 (For purposes of paragraph (1)(A)4. of this rule only, the term isomer includes the optical, position, and geometric isomers.):

A. Alpha-ethyltryptamine 7249
Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethenamine; 3-(2-aminobutyl)indole; alpha-ET; and AET;
B. 4-bromo-2,5-dimethoxyamphetamine 7391
Some trade or other names: 4-bromo-2, 5-dimethoxy-a-methylphenethylamine; 4-bromo-2, 5-DMA;
C. 4-bromo-2,5-dimethoxyphenethylamine 7392
D. 2,5-dimethoxyamphetamine 7396
Some trade or other names: 2,5-dimethoxyamphetamine; 2,5-DMA;
E. 2,5-dimethoxy-4-ethylamphetamine
Some trade or other names: DOET

F. 2,5-dimethoxy-4-(n)-propylthiophenethylamine
(other name: 2C-T-7)

G. 2-(2,5-Dimethoxy-4-(n)-propylphenyl) ethanamine
(2C-P)

H. 2-(2,5-Dimethoxy-4-ethylphenyl) ethanamine
(2C-E)

I. 2-(2,5-Dimethoxy-4-methylphenyl) ethanamine
(2C-D)

J. 2-(2,5-Dimethoxy-4-nitrophenyl) ethanamine (2C-N)

K. 2-(2,5-Dimethoxyphenyl) ethanamine (2C-H)

L. 2-(4-Chloro-2,5-dimethoxyphenyl) ethanamine (2C-C)

M. 2-(4-Ethylthio-2,5-dimethoxyphenyl) ethanamine (2C-T-2)

N. 2-(4-Iodo-2,5-dimethoxyphenyl) ethanamine (2C-I)

O. 2-(4-Isopropylthio)-2,5-dimethoxyphenyl) ethanamine (2C-T-4)

P. 4-methoxyamphetamine

Some trade or other names: 4-methoxyamethylphenethylamine;
paramethoxyamphetamine; PMA;

Q. 5-methoxy-3,4-methylenedioxyamphetamine

R. 4-methyl-2,5-dimethoxyamphetamine

Some trade and other names: 4-methyl-2,5-dimethoxy-a-methylphenethylamine; DOM; and STP;

S. 3,4- methylenedioxyamphetamine

T. 3,4-methylenedioxyamphetamine(MDMA)

U. 3,4-methylenedioxy-N-ethylamphetamine
(also known as N-ethylalphamethyl-3,4 (methylenedioxy) phenethylamine,
V. N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxyalpha-methyl-3,4 (methylenedioxy) phenethylamine and N-hydroxy MDA) 7402
W. 3,4,5- trimethoxyamphetamine 7390
X. 5-MeO-DMT or 5-methoxy-N,N-dimethyltryptamine 7431
Y. Alpha-methyltryptamine 7432
Z. Bufotenine 7433

Some trade and other names: 3-(b-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;
AA. Diethyltryptamine 7434

Some trade and other names: N,N-Diethyltryptamine; DET;
BB. Dimethyltryptamine

Some trade or other names: DMT;
CC. 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeODIPT) 7439
DD. Ibogaine 7260

Some trade and other names: 7-Ethyl-6,6β,7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido [1′,2′:1,2] azepino [5,4-b] indole; Tabernanthe iboga;
EE. Lysergic acid diethylamide 7315
FF. Marihuana 7360

Some trade or other names: marijuana;
GG. Mescaline 7381
HH. Parahexyl 7374

Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl;
II. Peyote 7415

Meaning all parts of the plant presently classified botanically as Lophophora williamsii Lemaire, 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 such plant, its seeds or extracts;
JJ. N-ethyl-3-piperidyl
benzilate 7482
KK. N-methyl-3-piperidyl benzilate 7484
LL. Psilocybin 7437
MM. Psilocyn 7438
NN. Tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis 7370 plant), as well as synthetic equivalents of the substances contained in the cannabis plant or in the resinous extractives of such plant, and/or synthetic substances, derivatives and their isomers, or both, with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following:
   (I) 1 cis or trans tetrahydrocannabinol and their optical isomers;
   (II) 6 cis or trans tetrahydrocannabinol and their optical isomers;
   (III) 3,4 cis or trans tetrahydrocannabinol and its optical isomers; and
   (IV) Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions are covered.
OO. Ethylamine analog of phencyclidine 7455
Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine, N-(1-phenylcyclohexyl)-ethylamine, cyclohexamine, PCE;
   PP. Pyrrolidine analog of phencyclidine 7458
Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine PCPy, PHP;
   QQ. Thiophene analog of phencyclidine 7470
Some trade or other names: 1-(1-(2-thienyl)cyclohexyl)-piperidine, 2-thienyl analog of phencyclidine, TPCP, TCP;
   RR. 1-(1-(2-thienyl)cyclohexyl)pyrrolidine 7473
Some other names: TCPy.
SS. Salvia divinorum
TT. Salvinorin A
UU. 3-Fluoromethcathinone 1233
VV. 4-Fluoromethcathinone 1238
WW. Mephedrone, or 4-methylmethcathinone 1248
XX. Methylenedioxypyrovalerone, MDPV, or (1-(1,3-Benzodioxol-5-yl)-2-(1-pyrrolidinyl)-1-pentanone 7535
YY. Methylone, or 3,4-Methylenedioxymethcathinone 7540
ZZ. Quinolin-8-yl 1-pentyl-1Hindole-3-carboxylate (PB-22; QUPIC) 7222
AAA. Quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22) 7225
BBB. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (AB-FUBINACA) 7012
CCC. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (ADB-PINACA) 7035
DDD. (1-pentyl-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone (Other names: UR-144, 1-pentyl-3-(2,2,3,3-tetramethylcyclopropyl)indole) 7144
EEE. [1-(5-fluoro-pentyl)-1Hindol-3-yl](2,2,3,3-tetramethylcyclopropyl) methanone (Other names: 5-fluoro-UR-144, 5-F-UR-144, XLR11, 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropyl)indole) 7011
FFF. N-(1-adamantyl)-1-pentyl-1Hindazole-3-carboxamide (Other names: APINACA, AKB48) 7048
GGG. 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (Other names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5) 7538

HHH. 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82) 7537

III. 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36) 7536

JJJ. 4-methyl-N-ethylcathinone (Other names: 4-MEC; 2-(ethylamino)-1-(4-methylphenyl)propan-1-one) 1249

KKK. 4-methyl-alpha-pyrrolidinopropiophenone, (Other names: 4-MePPP; MePPP; 4-methyl-alpha-pyrrolidinopropiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)-propan-1-one) 7498

LLL. alpha-pyrrolidinopentiophenone (Other names: alpha-PVP; alpha-pyrrolidinovalerophenone; 1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one) 7545

MMM. Butylone (Other names: bk-MBDB; 1-(1,3-benzodioxol-5-y1)-2-(methylamino)butan-1-one) 7541

NNN. Pentedrone (Other names: alpha-methylaminovalerophenone; 2-(methylamino)-1-phenylpentan-1-one) 1246

OOO. Pentyline (Other names: bk-
MBDP; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one)  

PPP. Naphyrone  
(Other names: naphthylpyrovalerone; 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one)  

QQQ. \( \alpha \)-pyrrolidinobutiophenone  
(Other names: \( \alpha \)-PBP; 1-phenyl-2-(pyrrolidin-1-yl)butan-1-one)  

RRR. \( N \)-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1Hindazole-3-carboxamide  
(Other names: AB-CHMINACA)  

SSS. \( N \)-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1Hindazole-3-carboxamide  
(Other names: AB-PINACA)  

TTT. [1-(5-fluoropentyl)-1Hindazole-3-yl](naphthalen-1-yl)methanone  
(Other names: THJ-2201)  

UUU. \( N \)-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1Hindazole-3-carboxamide  
(Other names: MAB-CHMINACA; ADB-CHMINACA)  

VVV. methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate  
(Other names: 5F–ADB; 5F–MDMB–PINACA)  

WWW. methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate  
(Other names: 5F–
AMB) 7033
XXX. N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide
   (Other names: 5F–APINACA, 5F–AKB48) 7049
YYY. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide
   (Other names: ADB–FUBINACA) 7010
ZZZ. methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate
   (Other names: MDMB–CHMICA, MMB–CHMINACA) 7042
AAAA. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate
   (Other names: MDMB–FUBINACA) 7020
BBBB. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate
   (Other names: FUB–AMB, MMB–FUBINACA, AMB–FUBINACA) (7021)

[UU]CCCC. 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 1H-indol-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:

(a) AM2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole 7201
(b) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole 7118
(c) JWH-015, or 1-propyl-2-methyl-3-(1-naphthoyl)indole 7019
(d) JWH-018, or 1-pentyl-3-(1-naphthoyl)indole 7173
(e) JWH-019, or 1-hexyl-3-(1-naphthoyl)indole 7081
(f) JWH-073, or 1-butyl-3-(1-naphthoyl)indole 7122
(g) JWH-081, or 1-pentyl-3-(4-methoxy-1-naphthoyl)indole 7122
(h) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole 7122
(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 7200
(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 7398
(m) JWH-398, or 1-pentyl-3-(4-chloro-1-naphthoyl)indole 7398

(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-morpholino)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-morpholino)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-morpholino)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-chlorophenylacetyl)indole
(c) JWH-250, or 1-pentyl-3-(2-methoxyphenylacetyl)indole
(d) JWH-251, or 1-pentyl-3-(2-methylphenylacetyl)indole
(e) RCS-8, or 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole

(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-morpholino)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.

- JWH-201
- JWH-203
- JWH-250
- JWH-251
- RCS-8
ethyl group, whether or not substituted in the cyclohexyl ring to any extent. Including, but not limited to:

(a) CP 47,497 & 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 7; 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-methyl octan-2-yl) - 6 a , 7 , 1 0 , 1 0 a - 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

5. 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 oxybutryrate; 2010
B. Mecloqualone 2572
C. Methaqualone 2565

6. 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. Aminorex 1585
Some trade or other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine;
B. N-benzylpiperazine (some other names: BZP, 1-benzylpiperzaine) 7493
C. Cathinone (Some trade or other names: 2-amino-1-phenyl-1-propanone, alphaaminopropiophenone, 2-aminopropiophenone and norephedrone) 1235
D. Fenethylline 1503

[E. 3-Fluoromethcathinone 1233
F. 4-Fluoromethcathinone 1238
G. Mephedrone, or 4-methylmethcathinone 1248]

[H]E. Methcathinone 1237
Some trade or other names: 2-(methylamino)-propiophenone; alpha-(methylamino) propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-N-methylaminopropiophenone; monomethylpropion; ephedrine; N-methylcathinone; methylcathinine; AL-464; AL-422; AL-463 and URI 432;

[J]F. 4-methoxymethcathinone

[K. Methylenedioxypyrovalerone,
MDPV, or (1-(1,3-Benzodioxol-5-yl)-2-(1-pyrrolidinyl)-1-pentanone

L. Methylone, or 3,4-Methylenedioxymethcathinone

[H] 4-Methyl-alphapyrrolidinobutiophenone, or MPBP

[N] I. N-ethylamphetamine 1475

[J] N,N-dimethylamphetamine 1480

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

[P] Quinolin-8-yl 1-pentyl-1Hindole-3-carboxylate (PB-22; QUPIC) 7222

Q. Quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22) 7225

R. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1Hindazole-3-carboxamide (AB-FUBINACA) 7012

S. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (ADB-PINACA) 7035

7. 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. (1-pentyl-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: UR-144, 1-pentyl-3-(2,2,3,3-tetramethylcyclopropyl)indole) 7144

B. [1-(5-fluoro-pentyl)-1Hindol-3-yl](2,2,3,3-tetramethylcyclopropyl) methanone, its optical,
position, and geometric isomers, salts, and salts of isomers (Other names: 5-fluoro-UR-144, 5-F-UR-144, XLR11, 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole)  
C. N-(1-adamantyl)-1-pentyl-1Hindazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomer (Other names: APINACA, AKB48)  
D. 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5)  
E. 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82)  
F. 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36)  
G. 4-methyl-N-ethylcathinone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: 4-MEC; 2-(ethylamino)-1-(4-methylphenyl)propan-1-one)  
H. 4-methyl-alphapyrrolidinopropiophenone, its optical, positional, and geometric isomers, salts, and
salts of isomers (Other names: 4-MePPP; MePPP; 4-methyl-α-pyrrolidinopropiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)-propan-1-one)

I. alphapyrrolidinopentiophenone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: α-PVP; α-pyrrolidinovalerophenone; 1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one)

J. Butylone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: bk-MBDB; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one)

K. Pentedrone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: α-methylaminovalerophenone; 2-(methylamino)-1-phenylpentan-1-one)

L. Pentylone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: bk-MBDP; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one)

M. Naphyrone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: naphthylpyrovalerone; 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one)

N. alpha-pyrrolidinobutiophenone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names:
α-PBP; 1-phenyl-2-(pyrrolidin-1-yl)butan-1-one) 7546

O. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1Hindazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-CHMINACA) 7031

P. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1Hindazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: AB-PINACA) 7023

Q. [1-(5-fluoropentyl)-1Hindazol-3-yl](naphthalen-1-yl)methanone, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: THJ-2201) 7024

R. N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (Other names: beta-hydroxythiofentanyl) 9836

[S]A. N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: acetyl fentanyl) 9821

[T] A. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1Hindazole-3-carboxamide, its optical, positional, and geometric isomers, salts, and salts of isomers (Other names: MAB-CHMINACA; ADB-CHMINACA) 7032
U. methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5F–ADB; 5F–MDMB–PINACA) (7034)

V. methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5F–AMB) (7033)

W. N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5F–APINACA, 5F–AKB48) (7049)

X. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: ADB–FUBINACA) (7010)

Y. methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: MDMB–CHMICA, MMB–CHMINACA) (7042)

Z. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate, its optical, positional, and geometric
isomers, salts and salts of isomers (Other names: MDMB–FUBINACA) (7020)

AA. N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other names: orthofluorofentanyl, 2-fluorofentanyl) (9816)

BB. 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: methoxyacetyl fentanyl) (9825)

CC. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: FUB–AMB, MMB–FUBINACA, AMB–FUBINACA) (7021)

DD. N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: cyclopropyl fentanyl) (9845)

[EE] B. N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: valeryl fentanyl) [(9804)]

[FF. N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl) butyramide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers]
(Other name: para-fluorobutyryl fentanyl)

\[ GG \] C. N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: para-methoxybutyryl fentanyl) (9823)

\[ HH \] D. N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: para-chloroisobutyryl fentanyl) (9837)

\[ II \] E. N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: isobutyryl fentanyl) (9826)

\[ JJ \] F. N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other name: cyclopentyl fentanyl) (9847)

\[ KK \] G. Fentanyl-related substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. (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.

\[ LL \]
Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: NM2201; CBL2201)

\[ MM \]
N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5F-AB-PINACA)

\[ NN \]
1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 4-CN-CUMYLBUTINACA; 4-cyano-CUMYL-BUTINACA; 4-CN-CUMYLBINACA; CUMYL-4CNBINACA; SGT-78)

\[ OO \]
methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoate, its
optical, positional, and geometric isomers, salts and salts of isomers (Other names: MMB-CHMICA, AMB-CHMICA)

\( [PP] \) L. 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5F-CUMYL-P7AICA)

\( [QQ] \) M. N-Ethylpentylone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: ephylone, 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one)

N. ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (trivial name: 5F-EDMB-PINACA)

O. methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate, its optical, positional, and geometric isomers, salts and salts of isomers (trivial name: 5F-MDMB-PICA)

P. N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (trivial names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-FLUOROBENZYL))

Q. 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (trivial names: 5F-CUMYL-PINACA; SGT-25)

R. (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,
<table>
<thead>
<tr>
<th>Path</th>
<th>Description</th>
<th>Chemical Name</th>
<th>Other Names</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>S.</td>
<td>N-Ethylhexedrone, its optical, positional, and geometric isomers, salts and salts of isomers (Other name: 2-(ethylamino)-1-phenylhexan-1-one)</td>
<td>3-tetramethylcyclopropyl) methanone, its optical, positional, and geometric isomers, salts and salts of isomers (trivial name: FUB-144)</td>
<td>7014</td>
<td></td>
</tr>
<tr>
<td>T.</td>
<td>Alpha-Pyrrolidinoheptaphenone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: alpha-PHP; alpha-pyrrolidinoheptaphenone; 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one)</td>
<td>S. N-Ethylhexedrone, its optical, positional, and geometric isomers, salts and salts of isomers (Other name: 2-(ethylamino)-1-phenylhexan-1-one)</td>
<td>7246</td>
<td></td>
</tr>
<tr>
<td>U.</td>
<td>4-Methyl-alpha-ethylaminopentiophenone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 4-MEAP; 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one)</td>
<td>7245</td>
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</tr>
<tr>
<td>V.</td>
<td>4′-Methyl-alpha-pyrrolidinohexiophenone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: MPHP; 4′-methyl-alpha-pyrrolidinohexiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one)</td>
<td>7446</td>
<td></td>
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</tr>
<tr>
<td>W.</td>
<td>Alpha-Pyrrolidinoheptaphenone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: PV8; 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one)</td>
<td>7446</td>
<td></td>
<td></td>
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<tr>
<td>X.</td>
<td>4-Chloro-alpha-pyrrolidinovalerophenone, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 4-chloro-alpha-PVP; 4′-chloro-alpha-pyrrolidinopentiophenone; 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one)</td>
<td>7548</td>
<td></td>
<td></td>
</tr>
<tr>
<td>Y.</td>
<td>N,N-diethyl-2-(2-(4 isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other names: isotonitazene; N,N-diethyl-2-[(4- (1-methylethoxy)phenyl)methyl]- 5-nitro-1H-benzimidazole-1- ethanamine)</td>
<td>9614</td>
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</tbody>
</table>
8. 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.

(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 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
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

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)

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, dextrophan, and levoproxyphene excepted:

A. Alfentanil
B. Alphaprodine
C. Anileridine
D. Bezitramide
E. Bulk Dextropropoxyphene (Non-dosage Forms)
F. Carfentanil
G. Dihydrocodeine
<p>| | |</p>
<table>
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<tbody>
<tr>
<td>A. Amphetamine, its salts, optical isomers, and salts of its optical isomers</td>
<td>1100</td>
</tr>
<tr>
<td>B. Lisdexamfetamine, its salts,</td>
<td></td>
</tr>
</tbody>
</table>

### 3. 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**
- **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:

A. Amobarbital 2125
B. Glutethimide 2550
C. Pentobarbital 2270
D. Phencyclidine 7471
E. Secobarbital 2315

5. Hallucinogenic substances:

A. Nabilone 7379
Another name for nabilone: (±)trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1-hydroxy-6,6-dimethyl-9H-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 and methamphetamine:
   (I) Phenylacetone 8501
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-piperidinocyclohexane carbonitrile (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 section 308.32 and any other drug of the quantitative 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
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
   (II) Secobarbital
   (III) Pentobarbital
   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
   (II) Secobarbital
   (III) Pentobarbital
   or any salt of any of these drugs and approved by 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

E. Embutramide

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)

H. Lysergic acid

I. Lysergic acid amide

J. Methyprylon

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

L. Sulfondiethylmethane
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, flupyrazapnon.

3. Nalorphine 9400

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 1.8 grams 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 1.8 grams 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 9804
   C. Not more than 1.8 grams 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 g) or not more than twenty-five milligrams (25 mg) per dosage unit, with one (1) or more active nonnarcotic ingredients in recognized therapeutic amounts.

F. Not more than fifty milligrams (50 mg) of morphine per one hundred milliliters (100 mL) or per one hundred grams (100 g), with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts.

5. Any material, compound, mixture, or preparation containing any of the following narcotic drugs or their salts, as set forth below:

   A. Buprenorphine

6. Anabolic steroids. Unless specially excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation. DEA has assigned code 4000 for all anabolic steroids. Anabolic steroids. Any drug or hormonal substance, chemically and pharmacologically related to testosterone (other than estrogens, progestins, corticosteroids, and dehydroepiandrosterone) that promotes muscle growth, except an anabolic steroid which is expressly intended for administration through implants to cattle or other nonhuman species and which has been approved by the Secretary of Health and Human Services for that administration. If any person prescribes, dispenses, or distributes such steroid for human use, such person shall be considered to have prescribed, dispensed, or distributed an anabolic steroid within the meaning of this subdivision. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, esters, and ethers:

   A. 3β,17β-dihydroxy-5α-androstane
   B. 3α,17β-dihydroxy-5α-androstane
<table>
<thead>
<tr>
<th>Letter</th>
<th>Name</th>
</tr>
</thead>
<tbody>
<tr>
<td>C.</td>
<td>5α-androstan-3,17-dione</td>
</tr>
<tr>
<td>D.</td>
<td>1-androstenediol (3β,17β-dihydroxy-5α-androst-1-ene)</td>
</tr>
<tr>
<td>E.</td>
<td>1-androstenediol (3α,17β-dihydroxy-5α-androst-1-ene)</td>
</tr>
<tr>
<td>F.</td>
<td>4-androstenediol (3β,17β-dihydroxy-androst-4-ene)</td>
</tr>
<tr>
<td>G.</td>
<td>5-androstenediol (3β,17β-dihydroxy-androst-5-ene)</td>
</tr>
<tr>
<td>H.</td>
<td>1-androstenedione ([5α]-androst-1-en-3,17-dione)</td>
</tr>
<tr>
<td>I.</td>
<td>4-androstenedione (androst-4-en-3,17-dione)</td>
</tr>
<tr>
<td>J.</td>
<td>5-androstenedione (androst-5-en-3,17-dione)</td>
</tr>
<tr>
<td>K.</td>
<td>Bolasterone (7α,17α-dimethyl-17β-hydroxyandrost-4-en-3-one)</td>
</tr>
<tr>
<td>L.</td>
<td>Boldenone (17β-hydroxyandrost-1,4-diene-3-one)</td>
</tr>
<tr>
<td>M.</td>
<td>Boldione (androsta-1,4-diene-3,17-dione)</td>
</tr>
<tr>
<td>N.</td>
<td>Calusterone (7β,17α-dimethyl-17β-hydroxyandrost-4-en-3-one)</td>
</tr>
<tr>
<td>O.</td>
<td>Clostebol (4-chloro-17β-hydroxyandrost-4-en-3-one)</td>
</tr>
<tr>
<td>P.</td>
<td>Dehydrochloromethyltestosterone (4-chloro-17β-hydroxy-17α-methyl-androst-1,4-dien-3-one)</td>
</tr>
<tr>
<td>Q.</td>
<td>Desoxymethyltestosterone (17α-methyl-5α-androst-2-en-17β-ol) (a.k.a. madol)</td>
</tr>
<tr>
<td>R.</td>
<td>Δ1-dihydrotestosterone (a.k.a.'1-testosterone') (17β-hydroxy-5α-androst-1-en-3-one)</td>
</tr>
<tr>
<td>S.</td>
<td>4-dihydrotestosterone (17β-hydroxy-5α-androst-1-en-3-one)</td>
</tr>
<tr>
<td>T.</td>
<td>Drostanolone(17β-hydroxy-2α-methyl-5α-androstan-3-one)</td>
</tr>
<tr>
<td>U.</td>
<td>Ethylestrenol(17α-ethyl-17β-hydroxyestr-4-ene)</td>
</tr>
<tr>
<td>V.</td>
<td>Fluoxymesterone(9-fluoro-17α-methyl-11β,17β-dihydroxyandrost-4-en-3-one)</td>
</tr>
<tr>
<td>W.</td>
<td>Formebulone (Formebolone) (2-</td>
</tr>
</tbody>
</table>
formyl-17α-methyl-11α,17β-dihydroxyandrost-1,4-dien-3-one

\[W\] X. Furazabol (17α-methyl-17β-hydroxyandrostano[2,3-c]-furazan)

\[X\] Y. 13β-ethyl-17β-hydroxygon-4-en-3-one

\[Y\] Z. 4-hydroxytestosterone (4,17β-dihydroxyandrost-4-en-3-one)

\[Z\] AA. 4-hydroxy-19-nortestosterone (4,17β-dihydroxy-estr-4-en-3-one)

\[AA\] BB. Mestanolone (17α-methyl-17β-hydroxy-5α-androstan-3-one)

\[BB\] CC. Mesterolone (1α-methyl-17β-hydroxy-[5α]-androstan-3-one)

\[CC\] DD. Methandienone (17α-methyl-17β-hydroxyandrost-1,4-dien-3-one)

\[DD\] EE. Methandriol (17α-methyl-3β,17β-dihydroxyandrost-5-ene)

\[EE\] FF. Methasterone (2α,17α-dimethyl-5α-androstan-17β-ol-3-one)

\[FF\] GG. Methenolone (1α-methyl-17β-hydroxy-5α-androst-1-en-3-one)

\[GG\] HH. 17α-methyl-3β,17β-dihydroxy-5α-androstan-4-ene

\[HH\] II. 17α-methyl-3α,17β-dihydroxy-5α-androstan-4-ene

\[II\] JJ. 17α-methyl-3β,17β-dihydroxyandrost-4-ene

\[JJ\] KK. 17α-methyl-4-hydroxynandrolone (17α-methyl-4-hydroxy-17β-hydroxyestr-4-en-3-one)

\[KK\] LL. Methyldienolone (17α-methyl-17β-hydroxyestr-4,9(10)-dien-3-one)

\[LL\] MM. Methyltrienolone (17α-methyl-17β-hydroxyestr-4,9,11-trien-3-one)

\[MM\] NN. Methyltestosterone (17α-methyl-17β-hydroxyandrost-4-en-3-one)

\[NN\] OO. Mibolerone (7α,17α-dimethyl-17β-hydroxyestr-4-en-3-one)

\[OO\] PP. 17α-methyl-D1 dihydrotestosterone (17β-hydroxy-17α-methyl-5α-androst-1-en-3-one) (a.k.a. 17α-methyl-1-testosterone)

\[PP\] QQ. Nandrolone (17β-hydroxyestr-4-ene-3-one)

\[QQ\] RR. 19-nor-4-androstenediol (3β,17β-dihydroxyestr-4-ene)
19-nor-4-androstenediol (3α,17β-dihydroxyestr-4-ene)
19-nor-4,9(10)-androstadienedione (estr-4,9(10)-dien-3,17-dione)
19-nor-5-androstenediol (3β,17β-dihydroxyestr-5-ene)
19-nor-5-androstenediol (3α,17β-dihydroxyestr-5-ene)
19-nor-4-androstenedione (estr-4-en-3,17-dione)
19-nor-5-androstenedione (estr-5-en-3,17-dione)
Norbolethone (13β,17α-diethyl-17β-hydroxygon-4-en-3-one)
Norclostebol (4-chloro-17β-hydroxyestr-4-en-3-one)
Norethandrolone (17α-ethyl-17β-hydroxyestr-4-en-3-one)
Normethandrolone (17α-methyl-17β-hydroxyestr-4-en-3-one)
Oxandrolone (17α-methyl-17β-hydroxy-2-oxa-[5α]-androst-1-en-3-one)
Oxymesterone (17α-methyl-4,17β-dihydroxyandrost-4-en-3-one)
Oxymetholone (17α-methyl-2-hydroxymethylene-17β-hydroxy-[5α]-androstan-3-one)
Prostanozol (17β-hydroxy-5α-androstan[3,2-c]pyrazole)
Stanolone (Δ1-dihydrotestosterone (a.k.a. 1-testosterone)(17β-hydroxy-5α-androst-1-en-3-one))
Stanozolol (17α-methyl-17β-hydroxy-[5α]-androst-2-eno[3,2-c]-pyrazole)
Stenbolone (17β-hydroxy-2-methyl-[5α]-androst-1-en-3-one)
Testolactone(13-hydroxy-3-oxo-13,17-secoandrost-1,4-dien-17-oic acid lactone)
Testosterone(17β-hydroxyandrost-4-en-3-one);
Tetrahydrogestrinone (13β,17α-diethyl-17β-hydroxygon-4,9,11-trien-3-one)
Trenbolone (17β-hydroxyestr-4,9,11-trien-3-one)
Any salt, ester, or isomer of a
drug or substance described or listed in this
subparagraph, if that salt, ester, or isomer
promotes muscle growth except an anabolic
steroid which is expressly intended for
administration through implants to cattle or
other nonhuman species and which has been
approved by the secretary of Health and
Human Services for that administration.

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

(Some other names for dronabinol: (6aRtrans)- 6a,7,8,10a-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 section. 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

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

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

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 g);

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

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

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:

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. Cloxazolam 2752
/O. Cloxazolam 2753
/P. Delorazepam 2754
/Q. Diazepam 2765
/R. Dichloralphenazone 2467
/S. Estazolam 2756
/T. Ethchlorvynol 2540
3. Fenfluramine. Any material, compound, mixture, or preparation which contains any quantity of the following substances, 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:

A. Fenfluramine 1670
4. 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

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. 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. Sibutramine 1675
   L. Solriamfetol (2-amino-3-phenylpropyl carbamate; benzenepropanol, beta-amino-, carbamate (ester)) 1650
   [L]M. SPA (-)-1-dimethylamino-1,2-diphenylethane 1635

6. 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 9725

7. 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 g);

B. Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 g);

C. Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 g);
D. Not more than two and five-tenths milligrams (2.5 mg) of diphenoxylate and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit;

E. Not more than one hundred milligrams (100 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 g); and

F. 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

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;

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:

A. Ezogabine [N-[2-amino-4(4-
fluorobenzylamino)-phenyl]-carbamic acid ethyl ester] 2779
B. Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide] 2746
C. Pregabalin [(S)-3-(aminomethyl)-5-methyl hexanoic acid] 2782
D. Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (also referred to as BRV; UCB-34714; Briviact) 2710
E. Lasmiditan [2,4,6-trifluoro-N-(6-(1-methylpiperidine-4-carbonyl)pyridine-2-yl-benzamide] 2790
F. 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
[5. Approved cannabidiol drugs.
A. A drug product in finished dosage formulation that has been approved by the U.S. Food and Drug Administration that contains cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-1,3-benzenediol) derived from cannabis and no more than one tenth percent (0.1%) (w/w) residual tetrahydro cannabinoids 7367]


PUBLIC COST: This emergency amendment will not cost state agencies or political subdivisions more than five hundred dollars ($500) annually.

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