# ARKANSAS REGISTER



# **Proposed Rule Cover Sheet**

Secretary of State John Thurston 500 Woodlane Street, Suite 026 Little Rock, Arkansas 72201-1094 (501) 682-5070 www.sos.arkansas.gov



Name of Department
Agency or Division Name
Other Subdivision or Department, If Applicable
Previous Agency Name, If Applicable
Contact Person_
Contact E-mail
Contact Phone_
Name of Rule
Newspaper Name
Date of Publishing
Final Date for Public Comment
Location and Time of Public Meeting

#### NOTICE OF PUBLIC HEARING

The Arkansas Department of Health will hold a public hearing at 1:30 pm on January 7, 2025, at the Arkansas Department of Health, Room 2508, located at 4815 West Markham, Little Rock, Arkansas, 72205 to allow interested persons to comment on the proposed amendments to the Rules Pertaining to the List of Controlled Substances in Arkansas.

Copies of the proposed Rules are available at <a href="www.healthy.arkansas.gov">www.healthy.arkansas.gov</a> and at Pharmacy Services, Room #357, 3<sup>rd</sup> Floor South, at the Arkansas Department of Health, 4815 W. Markham St. Little Rock, AR 72205.

The public may submit written comments no later than 4:30 p.m. on January 7, 2025. Comments may be sent by email to <a href="mailto:shane.david@arkansas.gov">shane.david@arkansas.gov</a> or by mail to Shane David, PharmD, Pharmacy Services Section, Arkansas Department of Health, 4815 West Markham Street, Slot #25, Little Rock, Arkansas, 72205.

List Of Controlled Substances



For the State Of Arkansas Pursuant to the provisions of Arkansas Code Annotated § 5-64-201 and § 5-64-216 of the laws of the State of Arkansas, the Secretary of the Arkansas Department of Health or duly authorized agent, as specified by law, is giving public notice of the publication of the List of Controlled Substances for the State of Arkansas.

Due consideration has been given applicable federal regulations, current scientific knowledge regarding the listed substances, the evidence of actual or relative potential for abuse, the history and current patterns of abuse, the risk to the public health, and potential to produce psychic or psychological dependence liability.

Based on these considerations the attached listing of the Schedule of Controlled Substances and the corresponding drugs that are included in each schedule is hereby promulgated by the Secretary of the Arkansas Department of Health as the List of Controlled Substances for the State of Arkansas.

Each controlled substance or basic class thereof has been assigned an "Administration Controlled Substance Code Number" for purposes of identification. These numbers are for internal management and are used as a means to identify substances with complex and cumbersome chemical names.

Next to the code number is the date the substance was placed in schedule by the Secretary of the Arkansas Department of Health. A "\*" denotes the substance was scheduled prior to April, 1979.

I, Shane David, Pharm.D., <u>Branch Chief, Health Systems Licensing and Certifications</u> <u>Section Chief of Pharmacy Services</u> for the Arkansas Department of Health, do hereby certify that the documents attached hereto are true and correct copies of the current List of Controlled Substances adopted by the Arkansas State Board of Health in accordance with Arkansas state law.

	Shane David, Pharm.D., Branch Chief Pharmacy Services Section
STATE OF ARKANSAS ) COUNTY OF SALINE )	
before me and signed the above	ereby certify that Shane David, Pharm.D., well known to me, appeared referenced document.  before me this day of March, 2024.
My commission expires	Notary Public

#### ARKANSAS DEPARTMENT OF HEALTH

#### LIST OF CONTROLLED SUBSTANCES

#### **SECTION I AUTHORITY**

The following scheduling of these controlled substances has been hereby promulgated pursuant to Arkansas Code Annotated §5-64-201 and §5-64-216.

#### SECTION II PURPOSE

Due consideration has been given applicable Federal regulations, current scientific knowledge regarding the listed substances, the evidence of actual or relative potential for abuse, the history and current patterns of abuse, the risk to the public health, and potential to produce psychic or psychological dependence liability.

## SECTION III GENERAL REQUIREMENTS

(Attached copy of the listing of scheduling of controlled substances)

# **SECTION IV REPEAL**

All lists of schedules of controlled substances in conflict herewith are hereby repealed.

#### **CERTIFICATION**

This will certify the following list of scheduling of controlled substances was adopted by the Arkansas State Board of Health at a session of the Board held in Little Rock, Arkansas on the 26<sup>th</sup> day of October, 20232024, and after a Public Hearing on the 23rd day of January, 2024, held in Little Rock, Arkansas, at the State Department of Health Building.

Jennifer Dillaha, MD Secretary of Arkansas State Board of Health Director of the Arkansas Department of Health

#### **ARTICLE II**

# **SCHEDULE I**

- (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 DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Opiates: (Narcotic Drugs)</u> 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, salts is possible within the specific chemical designation (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

(1) A	cetylmethadol	9601*
(2) A	llylprodine	9602*
(3) A	lphacetylmethadol (except Levo-alphacetylmethadol (LAAM)	-9603*
(4) A	lphameprodine	9604*
(5) A	lphamethadol	9605*
(6) B	enzethidine	9606*
(7) B	etacetylmethadol	9607*
(8) B	etameprodine	9608*
(9) B	etamethadol	9609*
(10)	Betaprodine	9611*
(11)	Dextromoramide	9613*
(12)	Diampromide	9615*
(13)	Diethylthiambutene	9616*
(14)	Difenoxin	9168*
(15)	Dimenoxadol	9617*
(16)	Dimepheptanol	9618*
(17)	Dimethylthiambutene	9619*
(18)	Dioxaphetyl butyrate	9621*
(19)	Dipipanone	9622*
(20)	Ethylmethylthiambutene	9623*
(21)	Etoxeridine	9625*
(22)	Furethidine	9626*
(23)	Hydroxypethidine	9627*

(24)	Ketobemidone	9628*
(25)	Levomoramide	9629*
(26)	Levophenacylmorphan	9631*
(27)	Morpheridine	9632*
(28)	MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)]	9661-(10-1985)
(29)	Noracymethadol	9633*
(30)	Norlevorphanol	9634*
(31)	Normethadone	9635*
(32)	Norpipanone	9636*
(33)	PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] - 9	9663-(10-1985)
(34)	Phenadoxone	9637*
(35)	Phenampromide	9638*
(36)	Phenomorphan	9647*
(37)	Phenoperidine	9641*
(38)	Piritramide	9642*
(39)	Proheptazine	9643*
(40)	Properidine	9644*
(41)	Propiram	9649*
(42)	Racemoramide	9645*
(43)	Tilidine	9750-(9-1981)
(44)	Trimeperidine	9646*
(45)	Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide]	(4-2017)
(46)	AH-7921 [other name(s): 3,4-dichloro-N-[(1 <del>N-[(1-</del> dimethylamino)cyclohexylmethyl]benzamide]	9551-(4-2017)
	W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4-chlorophenyl)sulfonamide]	(4-2017)
	W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4-chlorophenyl)Sulfonamide]	(4-2017)
(49)	MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine]	9560-(4-2017)
(50) N	U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)-N-methylbenzamide]	9547-(4-2017)
	Fentanyl-related substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Fentanyl-related substance means any substance not otherwise listed, and for which no exemption or approval is in effect under section	

	of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that is turally related to fentanyl by one or more of the following modifications:	(4-2021)
(i)	Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;	
(ii)	Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iii)	Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iv)	Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; or	
(v)	Replacement of the N-propionyl group by another acyl group.	
(vi)	Fentanyl-related substances shall include, but are not limited to:	
(.	A) Acetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	9815-(2-1986)
(1	B) Alpha-methylfentanyl [other name(s): (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propronanilide; 1-(1-methyl-2-phenylethyl)-4_(N-propanilido)piperidine)]	- 9814-(6-1982)
(	C) Alpha-methylthiofentanyl [other name(s): N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide]	
(]	D) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	- 9830-(2-1986)
(1	E) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide]	- 9831-(2-1986)
()	F) 3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-piperidyl]-N-Phenylpropanamide)]	9813-(10-1985)
(	G) 3-methylthiofentanyl [other name(s): N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide]	- 9833-(2-1986)
(1	H) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide]	9812-(11-1986)
(.	[) Thiofentanyl [other name(s): (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide]	9835-(2-1986)
(.	J) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9821-(4-2017)
(1	K) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	9822-(4-2017)
()	L) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2 yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	-

(M) Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	(4-2017)
(N) Valeryl fentanyl [other name(s): N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide]	-9840-(4-2017)
(O) Furanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran-2-carboxamide]	-9834-(4-2017)
(P) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	-9827-(4-2017)
(Q) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	-9838-(4-2017)
(R) 4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	(4-2017)
(S) Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide]	-9823-(4-2017)
(T) Acryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]	9811-(6-2020)
(U) 4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9824-(6-2020)
(V) Tetrahydrofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide]	
(W) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide]	9845-(4-2021)
(X) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9825-(4-2021)
(Y) Ortho-fluorofentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9816-(4-2021)
(Z) Crotonyl fentanyl [other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide]	9844-(5-2022)
(AA) Cyclopentyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide]	9847-(5-2022)
(BB) Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9826-(5-2022)
(CC) Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	9837-(5-2022)
(DD) Beta-methyl fentanyl [other name(s): N-phenyl-N-(1-(2-phenylpropyl) piperidin-4-yl)propionamide]	9856-(5-2022)
(EE) Beta'-phenyl fentanyl [other name: N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide]	

(FF) 2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	- 9855-(5-2022)
(GG) 4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) piperidin-4-yl)-N-phenylacetamide]	
(HH) Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	
(II) Ortho-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9848-(5-2022)
(JJ)Ortho-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	- 9820-(5-2022)
(KK) Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	- 9817-(5-2022)
(LL) Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide]	- 9841-(5-2022)
(MM) Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl) N-phenylthiophene-2-carboxamide]	
(NN) Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4-yl)(phenyl)carbamate]	- 9851-(5-2022)
(OO) Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1 phenethylpiperidin-4-yl)acrylamide]	
(PP) Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	- 9853-(5-2022)
(QQ) Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	- 9854-(5-2022)
(RR) Meta-fluorofentanyl [other name(s): N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9857
(SS) Meta-fluoroisobutyryl fentanyl [other name(s): N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide	9858
(TT) Para-methoxyfuranyl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	9859
(UU) Para-methylcyclopropyl fentanyl [other name(s): N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide]	9865
(VV) 3-furanyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3-carboxamide]	
(WW) 2',5'-dimethoxyfentanyl [other name(s): N-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-N-phenylpropionamide]	9861
(XX) Isovaleryl fentanyl [other name(s): 3-methyl-N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide]	9862

	(YY) Ortho-fluorofuranyl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	<u>9863</u>
	(ZZ) Alpha'-methyl butyryl fentanyl [other name(s): 2-methyl-N-(1-	
	phenethylpiperidin-4-yl)-N-phenylbutanamide]	<u> 9864</u>
(52)	Zipeprol	9873
(53)	Brorphine	9098
s a	Benzimidazole-opioid substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Benzimidazole-opioid substances includes any substance, not otherwise listed or excepted, 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 structurally has benzimidazole ring with an ethylamine at its 1-position and benzyl group at its 2-position:	
(	i) With or without substitution on the benzimidazole;	
(	ii) With or without substitution at the ethylamine;	
(	iii) With or without inclusion of the ethylamine in a cyclic structure;	
(	iv) With or without substitution on the benzyl ring; or	
(	v) With or without replacement of the benzyl ring with an aromatic ring.	
(	vi) Benzimidazole-opioid substances shall include but are not limited to:	
	(A)4'-Hydroxy Nitazene;	
	(B) 5-Aminoisotonitazene;	
	(C) Butonitazene;	
	(D)Clonitazene;	9612 *
	(E) Etodesnitazene, [other name(s): Etazene];	<del> 9765</del>
	(F) Etonitazene;	9624*
	(G) Flunitazene;	
	(H) Isotonitazene;	9614
	(I) Isotodesnitazene;	
	(J) Metodesnitazene;	
	(K) Metonitazene;	9757
	(L) N-Desethyl Etonitazene;	
	(M) N-Desethyl Isotonitazene;	
	(N) N-Piperidinyl Etonitazene [other name(s): Etonitazepipne];	
	(O) N-Pyrrolidino Etonitazene [other name(s): Etonitazepyne];	9758
	(P) N-Pyrrolidino Protonitazene; and	
	(O) Protonitazene	9759

	y	1)piperazin-1-y1)butan-1-one]	9664
(c)	sched the ex	Im derivatives: (Narcotic Drugs) Unless specifically excepted or unless listed dule, any of the following opium derivatives, its salts, isomers, and salts of isomers is possible within the specification:	ers whenever
	(1) A	cetorphine	9319*
	(2) A	cetyldihydrocodeine	9051*
	(3) B	enzylmorphine	9052*
	(4) C	odeine methylbromide	9070*
	(5) C	odeine-N-Oxide	9053*
	(6) C	yprenorphine	9054*
	(7) D	Desomorphine	9055*
	(8) D	Pihydromorphine	9145*
	(9) D	Protebanol	9335*
	(10)	Etorphine (except hydrochloride salt)	9056*
	(11)	Heroin	9200*
	(12)	Hydromorphinol	9301*
	(13)	Methyldesorphine	9302*
	(14)	Methyldihydromorphine	9304*
	(15)	Morphine methylbromide	9305*
	(16)	Morphine methylsulfonate	9306*
	(17)	Morphine-N-Oxide	9307*
	(18)	Myrophine	9308*
	(19)	Nicocodeine	9309*
	(20)	Nicomorphine	9312*
	(21)	Normorphine	9313*
	(22)	Pholcodine	9314*
	(23)	Thebacon	9315*
	(24)	Mitragynine	(11-2015)
	(25)	7-Hydroxymitragynine	(11-2015)

(55) 2-Methyl AP–237 [other name(s): 1-(2-methyl-4-(3-phenylprop-2-en-1-

(d) **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 substance, 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 this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

(1) alpha-ethyltryptamine	7249-(12-1993)
Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-i ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET.	ndole-3-
(2) 4-bromo-2,5-dimethoxy-amphetamine	7391*
Some trade or other names: 4-bromo-2,5-dimethoxy-alphamethylphenethylamine; 4-bromo-2,5-DMA.	
(3) 4-bromo-2,5-dimethoxyphenethylamine	7392-(8-1995)
Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1 an alpha-desmethyl DOB; 2C-B, Nexus.	ninoethane;
(4) 2,5-dimethoxyamphetamine	7396*
Some trade or other names: 2,5-dimethoxy-alpha-methylphenethyla DMA.	mine; 2,5-
(5) 2,5-dimethoxy-4-ethylamphetamine	7399-(3-1988)
Some trade or other names: DOET.	
(6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine	7348-(1-2005)
Some trade or other names: 2C-T-7.	
(7) 4-methoxyamphetamine	7411*
Some trade or other names: 4-methoxy-alpha- methylphenethylami paramethoxyamphetamine; PMA.	ne;
(8) 5-methoxy-3,4-methylenedioxy-amphetamine	7401*
(9) 4-methyl-2,5-dimethoxyamphetamine	7395*
Some trade and other names: 4-methyl-2,5-dimethoxy-alphamethylphenethylamine; "DOM"; and "STP".	
(10) 3,4-methylenedioxy amphetamine	7400*
(11) 3,4-methylenedioxymethamphetamine	7405-(10-1985)
Some trade or other names: MDMA)	
(12) 3,4-methylenedioxy-N-ethylamphetamine	7404-(6-1990)
Some trade or other names: N- ethy-alpha-methyl-3,4 (methylenedic phenethylamine, N-ethyl MDA; MDE; MDEA.	oxy)
(13) N-hydroxy-3,4-methylenedioxyamphetamine	7402-(6-1990)
Some trade or other names: N-hydroxy-alpha-methyl-3,4(methylene phenethylamine; N-hydroxy MDA	edioxy)

(14)	3,4,5-trimethoxy amphetamine	7390*
(15)	5-methoxy-n,n-dimethyltryptamine 5-MeO-DMT	- 7431*(1-2011)
(16)	alpha-methyltryptamine	- 7432-(7-2005)
	Some trade or other names: AMT	
(17)	Bufotenine	7433*
	Some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine.	
(18)	Diethyltryptamine	7434*
	Some trade or other names: N,N-Diethyltryptamine;DET.	
(19)	Dimethyltryptamine	7435*
	Some trade or other names: DMT	
(20)	5-methoxy-N,N-diisopropyltryptamine	- 7439-(7-2005)
	Some trade or other names: 5-MeO-DIPT.	
(21)	Ibogaine	7260*
	Some trade and other names: 7-Ethyl-6,6 beta; 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.	
(22)	Lysergic acid diethylamide	7315*
(23)	Mescaline	7381*
(24)	Parahexyl	- 7374-(7-1983)
	Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl.	
(25)	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, salts, derivative, mixture or preparation of such plant, its seeds or extracts. (Interprets 21 USC 812 (c), Schedule I (c) (12)).	
(26)	N-ethyl-3-piperidyl benzilate	7482*
(27)	N-methyl-3-piperidyl benzilate	7484*
(28)	Psilocybin	7437*
(29)	Psilocyn	7438*
(30)	Ethylamine Analog of phencyclidine	7455*

	Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE.	
(31)	Pyrrolidine Analog of phencyclidine	7458*
	Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCPy; PHP	
(32)	Thiophene Analog of phencyclidine	7470*
	Some trade or other names: 1-[1-(2-thienyl) cyclohexyl] Piperidine; 2-Thienyl analog of phencyclidine; TPCP; TCP.	
(33)	1-[1-(2-Thienyl)cylcohexyl]pyrrolidine	- 7473-(9-1989)
	Some other trade or other names: TCPy.	
(34)	N,N-Diallyl-5-Methoxytryptamine;	(6-2012)
	Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT	
(35)	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 5C-NBOMe	
	2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5I-NBOMe	(8-2013)
(37)	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E	7509-(11-2013)
(38)	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D	7508-(11-2013)
(39)	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C	7519-(11-2013)
(40)	2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I	7518-(11-2013)
(41)	2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-2	7385-(11-2013)
(42)	2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-4	7532-(11-2013)
(43)	2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H	7517-(11-2013)
(44)	2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N	7521-(11-2013)
(45)	2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine; 2C-P	7524-(11-2013)
	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine; 25B-	(9-2018)
	2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25B-	(11-2018)
(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH	(11-2018)
	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5E-NBOMe	(7-2019)
	2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 25H-NBOMe	(7-2019)
(51)	2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]- phenol; 25C-	

	(52)	2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25H-NBOH	(7-2019)
	(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	1245-(5-2022)
		Some trade or other names: Para-methoxymethamphetamine; PMMA	
	(54)	2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one	7286
		Some trade or other names: Methoxetamine; MXE	
(e)	com depr	pressants: Unless specifically excepted or unless listed in another schedule, any napound, mixture, or preparation which contains any quantity of the following substressant effect on the central nervous system, including its salts, isomers, and salts enever the existence of such salts, isomers, and salts of isomers is possible within the mical designation:	tances having a of isomers
	i	gamma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate], and its known precursors and analogs. Precursors include but are not limited to: gamma-butyrolactone	
	(2)	Mecloqualone	2572*
	(3)		2565*
	1 1 1	Benzodiazepine substances, their isomers, esters, ethers, salts and salts of isomers esters and ethers. Benzodiazepine substances includes any substance, not otherwise listed or excepted, 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 structurally has a fused 1,4-diazepine and benzene ring structure with a phenyl connected to the diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or any combination thereof. Benzodiazepine substances shall include but are not limited to:	
	(	(i) Bromazolam;	
	(	(ii) Clonazolam;	(4-2021)
	(	(iii) Flualprazolam;	(4-2021)
	(	(iv) Flubromazepam;	
	(	(v) Flubromazolam;	(4-2021)
	(	(vi) Phenazepam;	(6-2012)
	(	(vii) Phenazolam [other name(s): Clobromazolam].	
	i 1	Thienodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Thienodiazepine substances includes any substance, not otherwise listed or excepted, 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 structurally has a fused 1,4-diazepine and thiophene ring structure with a phenyl connected to the 1,4-diazepine ring, with any substitution(s) or	

		phen	cement(s) on the 1,4-diazepine or thiophene ring, any substitution(s) on the yl ring, or any combination thereof. Thienodiazepine substances shall include re not limited to:	
		(i)	Etizolam	(12-2014)
(f)	Sti	mula	nts:	
	(1)	comp subst	ss specifically excepted or unless listed in another schedule, any material, bound, mixture, or preparation which contains any quantity of the following cances having a stimulant effect on the central nervous system, including its isomers, and salts of isomers:	
		(i)	Cathinone	1235-(3-1988)
		(ii)	$(\pm)$ CIS-4-Methylaminorex $[(\pm)$ CIS-4,5-dihydro-4- methyl-5-phenyl-2-	. = 0 / 5 / 0 0 0 0
		····	oxazolamine]	,
			Fenethylline	
		(iv)	N-Benzylpiperazine [other name(s): BZP, 1-Benzylpiperazine]	7493-(1-2005)
			Some trade or other name: BZP, 1-Benzylpiperazine	
		(v)	N-ethylamphetamine	1475-(6-1982)
		(vi)	N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers	9818-(2-1986)
		(vii)	N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropan-mide (thenylfentanyl), its optical isomers, salts and salts of isomers	9834-(2-1986)
		(viii)	N,N-Dimethylamphetamine [other name(s): N,N,Alpha-trimethylbenzeneethanamine; N,N,Alpha-trimethylphenethylamine], its salts, optical isomers, and salts of optical isomers	1480-(2-1989)
		(ix)	Methcathinone [other name(s):(some 2-Methylamine-Proprophenone, alpha (methylamino)- Proprophenone, 2 (methylamino)-1-phenylpropan-1-one, alpha-N-Methylaminopropiophenone, monomethylpropion, ephedrone, N-methylcathinone, methycathinone, AL-464, AL-422, AL-463 and UR-1431], its salts, optical isomers and salts of optical isomers	237-(12-1993)
		(x)	Aminorex [other name(s): aminoraphen, 2-amino-5 phenyl-2-oxazoline, or 4,5 dihyrdo-5-phenyl-2-oxazolamine], its salts, optical isomers, and salts of optical isomers	585-(12-1993)
		(xi)	4,4'-Dimethylaminorex — [other name(s): some 4,4'-DMAR, 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine, or 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine]	1595-(5-2022)
		(xii)	Amineptine	
		(xiii)	Mesocarb	1227
		(viv)	Methyl-N-ethylcathinone (MEC)	(6.2014)

(2) Any material, compound, mixture, or preparation, whether produced directly or indirectly from a substance of vegetable origin or independently by means of chemical synthesis or by a combination of extraction and chemical synthesis, that contains any quantity of the following substances, or that contains any of the following substances' analogs, salts, isomers, and salts of isomers when the existence of the analogs, salts, isomers, and salts of isomers when the existence of the analogs, salts, isomers, and salts of isomers is possible within the specific chemical designation, with the following chemical structure is included in Schedule I:  (i) 4-Methylmetheathinone (Mephedrone) 1248-(3-2011)  (ii) Methylenedioxypyrovalerone (MDPV) 3-4-Methylenedioxypyrovalerone (MDPV) 13-4-Methylenedioxypyrovalerone (MDPV) 15-4-Methylenedioxypyrovalerone (Methylone) 15-40-(3-2011)  (iv) 4-Methoxymetheathinone 16-4-Methylenedioxypyrovalerone (Methylone) 15-4-Methylenedioxypyrovalerone (Mpha-PVP) 15-4-Methylenedio		(xv)	Methiopropamine. [other name(s): N-methyl-1-(thiophen-2-yl)propan-2-amine]	1478
(ii) Methylenedioxypyrovalerone (MDPV)	(2)	indir chem conta follo exist speci	ectly from a substance of vegetable origin or independently by means of nical synthesis or by a combination of extraction and chemical synthesis, that aims any quantity of the following substances, or that contains any of the wing substances' analogs, salts, isomers, and salts of isomers when the ence of the analogs, salts, isomers, and salts of isomers is possible within the fic chemical designation, with the following chemical structure is included:	e
(iii) 3,4-Methylenedioxy-N-methylcathinone (Methylone)		(i)	4-Methylmethcathinone (Mephedrone)	- 1248-(3-2011)
(iv)       4-Methoxymethcathinone       (3-2011)         (v)       3-Fluoromethcathinone       (3-2011)         (vi)       4-Fluoromethcathinone       (3-2011)         (vii)       1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)       7541-(11-2014)         (viii)       Alpha-Pyrrolidinopentiophenone (Alpha-PVP)       7545-(11-2015)         (ix)       4-methyl-N-ethylcathinone (4-MEC)       1249-(9-2018)         (x)       4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)       7498-(9-2018)         (xi)       2-(methylamino)-1-phenylpentan-1-one (Pentedrone)       1246-(9-2018)         (xii)       1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)       7542-(9-2018)         (xiii)       4-fluoro-N-methylcathinone (4-FMC, Flephedrone)       1238-(9-2018)         (xiv)       3-fluoro-N-methylcathinone (3-FMC)       1233-(9-2018)         (xvi)       1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)       1258-(9-2018)         (xvi)       Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)       7546-(9-2018)         (xvii)3-methylmethcathinone (3-MMC)       1259         (xvii)(xviii)       2-compound, unless listed in another schedule or a legend drug, that is structurally derived from 2-Amino-phenyl-1-propanone by modification or by substitution:       (3-2012)         (A) In the phenyl ring to any ext		(ii)	Methylenedioxypyrovalerone (MDPV)	(3-2011)
(v)       3-Fluoromethcathinone       (3-2011)         (vi)       4-Fluoromethcathinone       (3-2011)         (vii)       1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)       7541-(11-2014)         (viii)       Alpha-Pyrrolidinopentiophenone (Alpha-PVP)       7545-(11-2015)         (ix)       4-methyl-N-ethylcathinone (4-MEC)       1249-(9-2018)         (x)       4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)       7498-(9-2018)         (xi)       2-(methylamino)-1-phenylpentan-1-one (Pentedrone)       1246-(9-2018)         (xii)       1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)       7542-(9-2018)         (xiii)       4-fluoro-N-methylcathinone (4-FMC, Flephedrone)       1238-(9-2018)         (xiv)       3-fluoro-N-methylcathinone (3-FMC)       1233-(9-2018)         (xv)       1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)       1258-(9-2018)         (xvi)       Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)       7546-(9-2018)         (xvii)3-methylmethcathinone (3-MMC)       1259         (xvii)(xviii)       2         compound, unless listed in another schedule or a legend drug, that is structurally derived from 2-Amino-phenyl-1-propanone by modification or by substitution:       (3-2012)         (A) In the phenyl ring to any extent with alkyl, alkoxy, alkylenedioxy, <td></td> <td>(iii)</td> <td>3,4-Methylenedioxy-N-methylcathinone (Methylone)</td> <td>- 7540-(3-2011)</td>		(iii)	3,4-Methylenedioxy-N-methylcathinone (Methylone)	- 7540-(3-2011)
(vi) 4-Fluoromethcathinone		(iv)	4-Methoxymethcathinone	(3-2011)
(vii)       1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)       7541-(11-2014)         (viii)       Alpha-Pyrrolidinopentiophenone (Alpha-PVP)       7545-(11-2015)         (ix)       4-methyl-N-ethylcathinone (4-MEC)       1249-(9-2018)         (x)       4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)       7498-(9-2018)         (xi)       2-(methylamino)-1-phenylpentan-1-one (Pentedrone)       1246-(9-2018)         (xii)       1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)       7542-(9-2018)         (xiii)       4-fluoro-N-methylcathinone (4-FMC, Flephedrone)       1238-(9-2018)         (xiv)       3-fluoro-N-methylcathinone (3-FMC)       1233-(9-2018)         (xv)       1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)       1258-(9-2018)         (xvi)       Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)       7546-(9-2018)         (xvii)3-methylmethcathinone (3-MMC)       1259         (xvii)(xviii)       A         compound, unless listed in another schedule or a legend drug, that is structurally derived from 2-Amino-phenyl-1-propanone by modification or by substitution:       (3-2012)         (A) In the phenyl ring to any extent with alkyl, alkoxy, alkylenedioxy,		(v)	3-Fluoromethcathinone	(3-2011)
(viii) Alpha-Pyrrolidinopentiophenone (Alpha-PVP)		(vi)	4-Fluoromethcathinone	(3-2011)
(ix) 4-methyl-N-ethylcathinone (4-MEC)		(vii)	1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)	7541-(11-2014)
(x) 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)		(viii)	Alpha-Pyrrolidinopentiophenone (Alpha-PVP)	7545-(11-2015)
(xii) 2-(methylamino)-1-phenylpentan-1-one (Pentedrone)		(ix)	4-methyl-N-ethylcathinone (4-MEC)	1249-(9-2018)
(xii) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)		(x)	4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)	7498-(9-2018)
(xiii) 4-fluoro-N-methylcathinone (4-FMC, Flephedrone)		(xi)	2-(methylamino)-1-phenylpentan-1-one (Pentedrone)	1246-(9-2018)
(xiv) 3-fluoro-N-methylcathinone (3-FMC)		(xii)	1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)	7542-(9-2018)
(xvi) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)		(xiii)	4-fluoro-N-methylcathinone (4-FMC, Flephedrone)	1238-(9-2018)
(xvi) Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)		(xiv)	3-fluoro-N-methylcathinone (3-FMC)	1233-(9-2018)
(xvii)3-methylmethcathinone (3–MMC)		(xv)	1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)	1258-(9-2018)
(xvii)(xviii)		(xvi)	Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)	7546-(9-2018)
(xvii)(xviii)		(xvii	)3-methylmethcathinone (3–MMC)	1259
structurally derived from 2-Amino-phenyl-1-propanone by modification or by substitution:(3-2012)  (A) In the phenyl ring to any extent with alkyl, alkoxy, alkylenedioxy,			)(xviii)	
			structurally derived from 2-Amino-phenyl-1-propanone by modification	
phenyl ring by one (1) or more other univalent substituents;		(.	haloalkyl or halide substituents, whether or not further substituted in the	
(B) At the 3-position with an alkyl substituent; or		(	B) At the 3-position with an alkyl substituent; or	

(C) At the nitrogen atom with alkyl or dialkyl groups, or by inclusion of the nitrogen atom in a cyclic structure.	
(xviii)(xix)	
(xviii)(xix)	7543-(7-2019)
<del>(xix)</del> (xx)	
-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)	7547-(4-2021)
<del>(xx)</del> (xxi)	
-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone)	7549-(4-2021)
(xxi)(xxii)	
-(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone; Alpha-Ethylaminohexanophenone)]	7246
(xxii)(xxiii)	
1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (Alpha-Pyrrolidinohexanophenone; Alpha-PHP)]	7544
(xxiii)(xxiv)	
2-(ethylamino)-1-(4-methylphenyl)pentan-1-one [other name(s): (4-Methyl-alpha-ethylaminopentiophenone; 4-MEAP)]	
<del>(xxiv)</del> (xxv)	
1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (4'-Methyl-alpha-pyrrolidinohexiophenone; MPHP)]	7446
(xxv)(xxvi)	
-phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha-Pyrrolidinoheptaphenone; PV8)]	7548
(xxvi)(xxvii)	
1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): (4'-Chloro-alpha-pyrrolidinovalerophenone; 4-chloro-alpha-PVP)]	7443
(xxvii)(xxviii)	
4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): alpha-PiHP, Alpha-Pyrrolidinoisohexanophenone]	

## **SCHEDULE II**

- (a) 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.
- (b) **Narcotic Drugs:** Substances, vegetable origin or chemical synthesis. Unless specifically excepted or unless listed in another schedule, 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 combination of extraction and chemical synthesis:
  - (1) Opium and opiate, and any salt, compound, derivative, or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrorphan,

nalbuphine, naldemedine, nalmefene, naloxegol, naloxone,  $6\beta$ -naltrexol, naltrexone and samidorphan, and their respective salts, but including the following:

	(i)	Raw opium	9600*
	(ii)	Opium extracts	9610*
	(iii)	Opium fluid extracts	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*(9-2007)
	(xvi)	Oxycodone	9143*
	(xvii)	Oxymorphone	9652*
	(xviii	) Thebaine	9333*
	(xix)	Tapentadol	9780-(5-2009)
	(xx)	Noroxymorphone	9668-(4-2021)
(2)	equiv	salt, compound, derivative, or preparation thereof which is chemically ralent or identical with any of the substances referred to in paragraph (b) (1) is section, except that these substances shall not include the isoquinoline bids of opium.*	
(3)	Opiu	m poppy and poppy straw.*	
(4)	leave deriv deriv	leaves (9040) and any salt, compound, derivative, or preparation of coca s, (including cocaine (9041) and ecgonine (9180) and their salts, isomers, atives and salts of isomers and derivatives), and any salt, compound, ative, or preparation thereof which is chemically equivalent or identical with 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;	*
	(ii)	[ <sup>123</sup> I]ioflupane; or	

	(i	ii) [ <sup>18</sup> F]FP–CIT.	
	SC	Concentrate or poppy straw (the crude extract of poppy straw in either liquid, olid or powder form which contains the phenanthrene alkaloids of the opium oppy),	9670.*
(c)	Opia follow when	tes: (Narcotic Drugs) Unless specifically excepted or unless in another schedule, are wing opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters are ever the existence of such isomers, esters, ethers, and salts is possible within the specifical designations:	ny of the
	(1) A	Alfentanil 973	7-(2-1987)
	(2) A	lphaprodine	9010*
	(3) A	nileridine	9020*
	(4) B	ezitramide	9800*
	(5) B	sulk Dextropropoxyphene (non-dosage forms) 927	3-(9-1981)
	(6) C	Carfentanil 974	3-(9-1988)
	(7) D	Pihydrocodeine	9120*
	(8) D	Piphenoxylate	9170*
	(9) Fe	entanyl	9801*
	(10)	Isomethadone	9226*
	(11)	Levo-alphacetylmethadol (LAAM) 9648	-(12-1993)
	(12)	Levomethorphan	9210*
	(13)	Levorphanol	9220*
	(14)	Metazocine	9240*
	(15)	Methadone	9250*
	(16)	Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane	9254*
	(17) ca	Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-arboxylic acid	9802*
	(18)	Pethidine (Meperidine)	9230*
	(19)	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine	9232*
	(20)	Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate	9233*
	(21)	Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid	9234*
	(22)	Phenazocine	9715*
	(23)	Piminodine	9730*
	(24)	Racemethorphan	9732*
	(25)	Racemorphan	9733*

	(26)	Remifentanil	9739-(11-1996)
	(27)	Sufentanil	9740-(9-1981)
	(28)	Thiafentanil	9729-(4-2021)
	(29)	Oliceridine	9245-(5-2022)
	(30)	Tianeptine	(5-2022)
(d)	comp	<b>ulants:</b> Unless specifically excepted or unless listed in another schedule, any mound, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system:	·
	(1) A	mphetamine, its salts, optical isomers, and salts of its optical isomers	1100*
	(2) M	lethamphetamine, its salts, isomers, and salts of its isomers	1105*
	(3) L	isdexamefetamine	1205*(7-2007)
	(4) Pl	henmetrazine and its salts	1631*
	(5) M	lethylphenidate	1724*
(e)	comp depre when	essants: Unless specifically excepted or unless listed in another schedule, any round, mixture, or preparation which contains any quantity of the following subsections sant effect on the central nervous system, including its salts, isomers, and salts ever the existence of such salts, isomers, and salts of isomers is possible within tical designation:	tances having a of isomers
	(1) A	mobarbital	2125*
	(2) G	lutethimide	2550-(2-1991)
	(3) Po	entobarbital	2270*
	(4) Pl	hencyclidine	7471*
	(5) So	ecobarbital	2315*
(f)	Hallu	icinogenic Substances:	
	(1) N	abilone	7379-(11-1987)
		[Other name(s) for nabilone: $(\pm)$ -trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one].	
	U	ronabinol in an oral solution in a drug product approved for marketing by the .S. Food and Drug Administration; [(-)-delta-9-trans-trahydrocannabinol(delta-9-THC)]	7365-(7-2019)
(g)	Imm	ediate Precursor: Unless specifically excepted or unless listed in another sched	lule, any

material, compound, mixture, or preparation which contains any quantity of the following

substances:

	(1) In	mmediate precursor to Amphetamine and Methamphetamine:	
	(i	i) Phenylacetone	8501-(3-1980)
		Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl Ketone; methyl benzyl Ketone.	
	(2) In	mmediate precursor to Phencyclidine (PCP):	
	(i	i) 1-phenylcyclohexylamine	····· 7460*
	(i	ii) 1-piperidinocyclohexanecarbonitrile (PCC)	8603*
	(3) In	mmediate precursor to Fentanyl:	
	(i	i) 4-anilino-N-phenethylpiperidine (ANPP) 8	3333*(8-2010)
	(i	ii) N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl)	8366-(4-2021)
SC	HED	<u>ULE III</u>	
(a)	usual	dule III shall consist of the drugs and other substances, by whatever official name, l name, chemical name, or brand name designated, listed in this section. Each drug een assigned the DEA Controlled Substances Code Number set forth opposite it.	
(b)	comp stimu or ge	cound, mixture, or preparation which contains any quantity of the following substantant effect on the central nervous system, including its salts, isomers (whether optermetric), and salts of such isomers whenever the existence of such salts, isomers, ers is possible within the specific chemical designation:	ances having a cical, position,
	st p S li	Those compounds, mixtures, or preparations in dosage unit form containing any timulant substances listed in Schedule II which compounds, mixtures, or reparations 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 ontrolled substances————————————————————————————————————	1405*
	(2) B	Benzphetamine	1228*
	(3) C	Chlorphentermine	1645*
	(4) C	Clortermine	1647*
	(5) P	PhendimetrazinePhendimetrazine	1615*
(c)	comp	ressants: Unless specifically excepted or unless listed in another schedule, any macound, mixture, or preparation which contains any quantity of the following substates and effect on the central nervous system:	
	(1) A	Any compound, mixture, or preparation containing:	
	(i	i) Amobarbital	2126*
	(i	ii) Secobarbital	2316*

(i	ii) Pentobarbital	2271*
(i	v) Embutramide	2020*(9-2006)
	or any salt thereof and one or more other active medicinal ingredients which are not listed in any schedule.	
(2) A	ny suppository dosage form containing:	
(i	) Amobarbital	2126*
(i	i) Secobarbital	2316*
(i	ii) Pentobarbital	2271*
	or any salt of any of these drugs and approved by the Food and Drug Administration for marketing only as a suppository.	
` '	ny substance which contains any quantity of a derivative of barbituric acid or	
	ny salt thereof	
. ,	hlorhexadol	2510*
` /	ny drug product containing gamma hydroxybutyric acid, including its salts, omers, and salts of isomers, for which an application is approved under section	
	Of the Federal Food, Drug, and Cosmetic Act	- 2012-(2-2001)
	etamine. its salts, isomers, and salts of isomers	
` ,	Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	, ,
(7) L	ysergic acid	7300*
(8) L	ysergic acid amide	7310*
(9) M	lethyprylon	2575*
(10)	Sulfondiethylmethane	2600*
(11)	Sulfonethylmethane	2605*
(12)	Sulfonmethane	2610*
(13)	Tiletamine and zolazepam or any salt thereof	
	Some trade or other name for a tiletamine- zolazepam 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,-trimethyl_pyrazolo-[3,4-e]-[1,4,]-diazepin-7(1-H)-one-; flupyrazapon.	
(14)	Perampanel	2261-(11-2013)
(15) <u>ar</u>	Xylazine and any material, compound, mixture, or preparation which contains by quantity of xylazine, including its salts, isomers, and salts of isomers	

(i) Dispensing, prescribing, or administering, to an animal, a drug containing xylazine that has been approved by the United States Secretary of Health and Human Services under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b);	
(ii) Dispensing, prescribing, or administering xylazine to an animal that is permissible under section 512 (a)(4) of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(a)(4));	
(iii) Possessing a drug containing xylazine, as described in this section (15), for animal use:	
(A) By a licensed pharmacist or licensed veterinarian; or	
(B) Pursuant to a valid prescription from a licensed veterinarian;	
(iv) Possessing, manufacturing, distributing, or using xylazine as an active pharmaceutical ingredient for manufacturing an animal drug either:	
(A) Approved under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b); or	
(B) Issued an investigation use exemption under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(j));	
(v) Manufacturing, distributing, or using a xylazine bulk chemical for pharmaceutical compounding by a licensed pharmacist or veterinarian; or	
(vi) Another use approved or permissible under the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 301, et seq.) or under 21 CFR Part 530, Subpart B.	
(d) Nalorphine	9400*
(e) Narcotic drugs: Unless specifically excepted or unless listed in another schedule:	
(1) Any material, 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 set forth below:	
(i) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*
(ii) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9804*
(iii) Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9807*

whenever the existence of such salts, isomers, and salts of isomers is possible

within the specific chemical designation, except in the following uses:

	(1)	more than 15 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9808*		
	(v	Not more than 500 milligrams of opium per 100 milliliters or per 100 grams or not more than 25 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts			
	(v	Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9810*		
		ny material, compound, mixture, or preparation containing any of the following rcotic drugs or their salts, as set forth below:			
	(i)	Buprenorphine (10-02 Transfer)	9064-(6-1985)		
	(ii	) Reserved			
(f)	composalts,	olic Steroids: Unless specifically excepted or unless listed in another schedule, a bund, mixture, or preparation containing any quantity of the following substance isomers, and salts of isomers whenever the existence of such salts of isomers is pecific chemical designation:	es, including its possible within		
	(1) Bo	oldenone;	(9-1991)		
	(2) Bo	oldione;	(1-2010)		
	(3) Cl	nlorotestosterone (4-chlortestosterone);	(9-1991)		
	(4) Clostebol;(9-199				
	(5) Dehydrochlormethyltestosterone;(9-1991				
	(6) Desoxymethyltestosterone(1-20				
	(7) D	hydrotestosterone (4-dihydrotestosterone);	(9-1991)		
	(8) D	rostanolone;	(9-1991)		
	(9) Et	hylestrenol;	(9-1991)		
	(10)	Fluoxymesterone;	(9-1991)		
	(11)	Formebulone (formebolone);	(9-1991)		
	(12)	Mesterolone;	(9-1991)		
	(13)	Methandienone, also known as Methandrostenolone;	(9-1991)		
	(14)	Methandranone;	(9-1991)		
	(15)	Methandriol;	(9-1991)		
	(16)	Methenolone;	(9-1991)		
	(17)	Methyltestosterone;	(9-1991)		

(18)	Mibolerone;	(9-1991)
(19)	Nandrolone;	(9-1991)
(20)	19-Nor-4,9(10)-Androstadienedione	(1-2010)
(21)	Norethandrolone;	(9-1991)
(22)	Oxandrolone;	(9-1991)
(23)	Oxymesterone;	(9-1991)
(24)	Oxymetholone;	(9-1991)
(25)	Stanolone;	(9-1991)
(26)	Stanozolol;	(9-1991)
(27)	Testolactone;	(9-1991)
(28)	Testosterone;	(9-1991)
(29)	Trenbolone	(9-1991)
(30)	Prostanozol	(8-2012)
(31)	Methasterone	(8-2012);
	and	
(g) Exen	aragraph, if that salt, ester, or isomer promotes muscle growth  ppt anabolic steroid products: Compounds, mixtures, or preparations that did that have been exempted by the Secretary:  NDC Number	
(1) <u>A</u>	ndro-Estro 90-4	0536-1605
` /	ndrogyn L.A	
. ,	omponent E-H in Process Pellets	
	omponent E-H in Process Granulation	-
` /	omponent TE-S in process Granulation	•
	omponent TE-S in process Pellets	•
	epANDROGYN	•
` '	epo-Testadiol	
` /	PEPO-T.E	
(10)	depTESTROGEN	
(11)	Duomone	
(12)	DUO-SPAN II	
(1-)	· · - · · · ·	000.0102
(13)	DURATESTRIN	43797-016

(14)	Essian	- Pharmaceutics
(15)	Essian H.S	- Pharmaceutics
(16)	Esterified Estrogens & Methyltestosterone, USP (0.625 mg/1.25mg)	Interpharm
(17)	Esterified Estrogens & Methyltestosterone, USP (1.25mg/2.5mg)	Interpharm
(18)	Esterified Estrogens & Methyltestosterone (0.625mg/1.25mg) Tablet	ANDAPharm
(19)	Esterified Estrogens & Methyltestosterone (1.25mg/2.5mg) Tablet	ANDAPharm
(20)	Estratest	0032-1026
(21)	Estratest HS	0032-1023
(22)	Menogen	59243-570
(23)	Menogen HS	59243-560
(24)	Methyltestosterone & Esterified Estrogens (2.5mg/1.25Mg)	Lannett Co
(25)	Methyltestosterone & Esterified Estrogens (Half Strength) (1.25mg/0.625mg)	Lannett Co
(26)	PAN ESTRA TEST	0525-0175
(27)	Premarin with Methyltestosterone	0046-0879
(28)	Premarin with Methyltestosterone	0046-0878
(29)	Syntest D.S	66576-231
(30)	Stntest H.S	66576-230
(31)	Synovex H in process bulk pellets	Syntex Animal
(32)	Synovex H in process granulation	Syntex Animal
(33)	Synovex Plus in process granulation	Fort Dodge
(34)	Synovex Plus in process bulk pellets	Fort Dodge
(35)	TEST-ESTRO Cypionates	0536-9470
(36)	Testoderm with Adhesive 4mg/d	Alza Corp
(37)	Testoderm 4mg/d	17314-4608
(38)	Testoderm 6mg/d	17314-4609
(39)	Testoderm with Adhesive 6mg/d	17314-2836
(40)	Testoderm in process film	Alza Corp
(41)	Testoderm with Adhesive in process film	Alza Corp
(42)	Testosterone Cypionate/Estradiol Cypionate injection	54274-530
(43)	Testosterone Cypionate/Estradiol Cypionate injection	0182-3069
(44)	Testosterone Cyp 50 Estradiol Cyp 2	0814-7737
(45)	Testosterone Cypionate/Estradiol Cypionate injection	0364-6611
(46)	Testosterone Cypionate/Estradiol Cypionate injection	0402-0257

(47) Testosterone Enanthate/Estradiol Valerate injection	0182-3073
(48) Testosterone Enanthate/Estradiol Valerate injection	0364-6618
(49) Testosterone Enanthate/Estradiol Valerate injection	0402-0360
(50) Testosterone Ophthalmic Solution	Allergan
(51) Tilapia Sex Reversal Feed (investigational)	Ranger, Inc
(h) Veterinary Anabolic Steroid Implant Products: Anabolic steroid products administration through implants in cattle or other nonhuman species exempted.	
NDC/DIN	
(1) Component E-H	021641-002
(2) Component E-H	01968327
(3) Component TE-S	021641-004
(4) Component T-H	0211641-006
(5) Component T-S	0211641-005
(6) F-TO	00093351
(7) Finaplix-H	12799-807-10
(8) Finaplix-S	12799-807-07
(9) Heifer-old	Boehringer
(10) Heifer-old	Ingelheim
(11) Heifer-old	Ivy Lab.
(12) Implus-H	0009-0434-01
(13) Implus-H	06-0434-01
01968327	
(14) Masculinizing Feed for Fish (Invesitigational)	Rangen,Inc.
(15) Revalor-G	12799-811
(16) Revalor-H	12799-810
(17) Revalor-S	12799-809
(18) Synovex H	0856-3901
(19) Synovex H	Syntex
(20) Synovex Plus	0856-3904
(21) Tilapia Sex Reversal Feed (investigational)	Zeigier Bros.

If veterinary products that are granted exempted status are subsequently distributed with the intent that they be used in humans, the distribution would be subject to the criminal sanctions of the CSA despite the drugs' exempted status.

# (i) Hallucinogenic substances:

(1) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a U.S. Food and Drug Administration approved drug product ----- 7369-(11-1987)

[Some other names for dronabinol: (6a R-trans)-6a,7,8, 10a-tetrahydro-6, 6, 9-trimethyl-3-pentyl-6H-dibenzo [b,d] phyran-1-ol, or (-)-delta 9-(trans)-tetrahydrocannabinol]

#### **SCHEDULE IV**

- (a) 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.
- (b) <u>Narcotic drugs</u>: Unless specifically excepted or unless listed in another schedule, any material, 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 set forth below:
  - (1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit------ 9167\*
  - (2) Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-propionoxybutane) ------ 9278-(11-1987)
- (c) <u>Depressants</u>: 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:

(13) Delorazepam ------ 2754-(1-1985)

(14)	Diazepam	2765*
(15)	Dichloralphenazone	2467-(10-2002)
(16)	Estazolam	2756-(1-1985)
(17)	Ethchlorvynol	2540*
(18)	Ethinamate	2545*
(19)	Ethyl loflazepate	2758-(1-1985)
(20)	Fludiazepam	2759-(1-1985)
(21)	Flunitrazepam	2763-(1-1985)
(22)	Flurazepam	2767*
(23)	Fospropofol	2138-(11-2009)
(24)	Halazepam	2762-(6-1982)
(25)	Haloxazolam	2771-(1-1985)
(26)	Ketazolam	2772-(1-1985)
(27)	Loprazolam	2773-(1-1985)
(28)	Lorazepam	2885*
(29)	Lormetazepam	2774-(1-1985)
(30)	Mebutamate	2800*
(31)	Medazepam	2836-(1-1985)
(32)	Meprobamate	2820*
(33)	Methohexital	2264*
(34)	Methylphenobarbital (mephorbarbital)	2250*
(35)	Midazolam	2884-(1-1985)
(36)	Nimetazepam	2837-(1-1985)
(37)	Nitrazepam	2834-(1-1985)
(38)	Nordiazepam	2838-(1-1985)
(39)	Oxazepam	2835*
(40)	Oxazolam	2839*
(41)	Paraldehyde	2585*
(42)	Petrichloral	2591*
(43)	Phenobarbital	2285*
(44)	Pinazepam	2883-(1-1985)
(45)	Prazepam	2764*
(46)	Quazepam	2881-(11-1986)

(47	) Temazepam	2925-(9-1981)
(48	) Tetrazepam	2886-(1-1985)
(49	) Triazolam	2887-(7-1983)
(50	) Zaleplon	2781-(9-1999)
(51	) Zolpidem	2783-(12-1993)
(52	) Zopiclone	2784-(1-2006)
(53	) Alfaxalone	2731-(2-2014)
(54	) Carisoprodol	8192-(4-1997)
(55	) Tramadol	9752-(8-2007)
(56	) Suvorexant	2223-(8-2014)
(57	) Brexanolone	2400-(4-2021)
(58	) Lemborexant	2245-(4-2021)
(59	) Remimazolam	2846-(5-2022)
(60	) Daridorexant	2410
<u>(61</u>	Zuranolone	<u> 2420</u>
or g	nulant effect on the central nervous system, including its salts, isomers (vgeometric), and salts of such isomers whenever the existence of such salts mers is possible within the specific chemical designation:	
(1)	Cathine ((+)-Norpseudoephedrine)Norpseudeophedrine	1230-(3-1988)
(2)	Diethylpropion	1610*
(3)	Fencamfamin	1760-(3-1988)
(4)	Fenproporex	· · · · · · · · · · · · · · · · · · ·
(5)	Lorcaserin	1575-(3-1988)
(6)		1575-(3-1988)
(7)	Mazindol	1575-(3-1988) 1625-(6-2013)
( )	Mazindol Mefenorex	1575-(3-1988) 1625-(6-2013) 1605-(6-1982)
` '		1575-(3-1988) 1625-(6-2013) 1605-(6-1982) 1580-(3-1988)
(8)	Mefenorex	1575-(3-1988) 1625-(6-2013) 1605-(6-1982) 1580-(3-1988) 1680-(1-1999)
(8)	Mefenorex  Modafinil  Pemoline (including organometallic complexes and chelates thereof)	1575-(3-1988) 1625-(6-2013) 1605-(6-1982) 1580-(3-1988) 1680-(1-1999) 1530*
(8) (9)	Mefenorex  Modafinil  Pemoline (including organometallic complexes and chelates thereof)  Phentermine	1575-(3-1988) 1625-(6-2013) 1605-(6-1982) 1580-(3-1988) 1680-(1-1999) 1530*
(8) (9) (10	Mefenorex  Modafinil  Pemoline (including organometallic complexes and chelates thereof)  Phentermine  Pipradrol	1575-(3-1988) 1625-(6-2013) 1605-(6-1982) 1580-(3-1988) 1680-(1-1999) 1530* 1640* 1750-(9-1981)

	(14) Solriamfetol	1650-(4-2021)
	(15) SPA ((-)-1-dimethylamino-1,2,diphenylethane)	1635-(9-1981
e)	Other substances: Unless specifically excepted or unless listed in another schedule, compound, mixture, or preparation which contains any quantity of the following substincluding its salts; isomers whether optical, position, or geometric), and salts of such its whenever the existence of such salts, isomers, and salts of isomers is possible:	tances,
	(1) Pentazocine	9709-(4-1979
	(2) Butorphanol	9720-(4-1997
	(3) Nalbuphine	(4-1997
	(4) Eluxadoline	9725-(4-2017
SC	CHEDULE V	
a)	Schedule V shall consist of the drugs and other substances by whatever official name, usual name, chemical name, or brand name designated, listed in this section.	common or
(b)	<b>Narcotic Drugs:</b> Unless specifically excepted or unless listed in another schedule, are compound, mixture or preparation containing any of the following narcotic drugs and set forth below.	•
	Reserved	
(c)	Narcotic drugs containing nonnarcotic active medicinal ingredients. Any composition or preparation containing any of the following limited quantities of narcotic drugs or such which shall include one or more nonnarcotic active medicinal ingredients in sufficient confer upon the compound, mixture, or preparation valuable medicinal qualities other possessed by the narcotic drug alone:	salts thereof, t proportion to
	(1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams	;
	(2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams	;
	(3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams.	;
	(4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit.	;
	(5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams	;
	(6) Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.	;
d)	<u>Stimulants</u> : unless specifically excepted or unless listed in another schedule, any ma compound, mixture, or preparation which contains any quantity of the following substimulant effect on the central nervous system, including its salts, isomers and salts of	tances having

(1) Pyrovalerone	1485-(3-1988)
(2) Ephedrine:a -{-(Methylamino)ethyl}benzene-methanol;	(10-1995)
a-{-(methylamino) ethyl} benzyl alcohol; 2-methylamino-1-phenyl-1-propanol; 1-phenyl-1-hydroxy-2-methylaminopropane; 1-phenyl-2-methylaminopropanol; a - hydroxy-b-methylaminopropylbenzene; a product which occurs in the Chinese herb Ma Huang (Ephedra vulgaris, Ephedra sinica Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra spp.	
(3) Phenylpropanolamine	(7-2005)
(4) Pseudoephedrine	(7-2005)
Pursuant to Ark. Code Ann. § 5-64-212 as amended in 2005, this Schedule V classific NOT apply to any ephedrine, phenylpropanolamine, or pseudoephedrine in liquid, liquid gel capsule form. However, sales limits mandated by statute shall apply to all pephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless form.	uid capsule, or products with
(e) <b>Depressants:</b> Unless specifically exempted or excluded or unless listed in another solution material, compound, mixture, or preparation which contains any quantity of the follow having a depressant effect on the central nervous system, including its salts, isomers, a isomers:	ving substances
(1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	2782-(1-2006)
(2) Lacosamide	2746-(5-2009)
(3) Brivaracetam	2710-(4-2017)
(4) Lasmiditan	2790-(4-2021)
(5) Cenobamate	2720-(4-2021)
(6) Ganaxolone	2401

## (f) Other substances:

(1) None.

# **SCHEDULE VI \*\*\*\***

(a) In addition to any substance placed in Schedule VI by the Secretary of the Department of Health under § 5-64-214, any material, compound, mixture, or preparation, whether produced directly or indirectly from a substance of vegetable origin or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis, that contains any quantity of the following substances, or that contains any of their salts, isomers, and salts of isomers when the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation, is included in Schedule VI:

1)!	Ma	arijuana	*	*
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(2)	Tetra	ahydro	ocannabinols, unless the tetrahydrocannabinol is:	**
	(i)	Cont	tained in hemp-derived cannabidiol;	(6-2020)
	(ii)	tet bas	more than three-tenths of one percent (0.3%) of delta-9 rahydrocannabinol in the hemp-derived cannabidiol on a dry weight sis as verified by a nationally accredited laboratory for quality, purity d accuracy standards; and	(6-2020) ***
	(iii)		approved by the United States Food and Drug Administration for urketing as a medication;	(6-2020)
(3)	A sy	ntheti	c equivalent of:	
	(i)	The	substance contained in the Cannabis plant; or	**
	(ii)	The	substance contained in the resinous extractives of the genus Cannabis;	**
(4)	class the p manu extra salts.	ified lant, a	botanically as Salvia divinorum, whether growing or not, the seeds of any extract from any part of the plant, and every compound, are, derivative, mixture, or preparation of the plant, its seeds, or its including salts, isomers, and salts of isomers when the existence of the ners, and salts of isomers is possible within the specific chemical	**
			n;	**
(3)	class specific Com nume The	es des ific un pound erical	substances, derivatives, or their isomers in the chemical structural scribed below in subdivisions $(a)(5)(i)-(a)(5)(x)$ of this section and also neclassified substances in subdivision $(a)(5)(xi)$ of this section. ds of the structures described in this subdivision $(a)(5)$ , regardless of designation of atomic positions, are included in this subdivision $(a)(5)$ . etic substances, derivatives, or their isomers included in this subdivision	
	(i)	Tetra	ahydrocannabinols:	
	(.	A)Te	trahydrocannabinols, including without limitation the following:	**
		a)	Delta-1 cis or trans tetrahydrocannabinol [other name(s): Delta-9 cis or trans tetrahydrocannabinol], and its optical isomers;	**
		b)	Delta-6 cis or trans tetrahydrocannabinol [other name(s): Delta-8 cis or trans tetrahydrocannabinol], and its optical isomers;	**
		c)	Delta- 3,4 cis or trans tetrahydrocannabinol [other name(s): Delta-6a,10a cis or trans tetrahydrocannabinol], and its optical isomers;	**
		d)	Delta-10 cis or trans tetrahydrocannabinol, and its optical isomers;	***
		e)	Delta-8 tetrahydrocannabinol acetate ester;	***
		f)	Delta-9 tetrahydrocannabinol acetate ester;	***
		g)	Delta-6a,10a, tetrahydrocannabinol acetate ester;	
		h)	Delta-10 tetrahydrocannabinol acetate ester; and,	***

		A product derived from industrial hemp that was produced as a result of a synthetic chemical process that converted the industrial hemp or a substance contained in industrial hemp into Delta-8, Delta-9, Delta 6a,10a, or Delta-10 tetrahydrocannabinol including their respective acetate esters.	***
		ronabinol in sesame oil and encapsulated in a soft gelatin capsule in a ug product approved by the United States Food and Drug dministration is not a tetrahydrocannabinol under this subdivision ()(5)(i);	**
(ii)		hthoylindoles, or any compound structurally derived from 3-(1-phthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at enitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-morpholinyl)ethyl group, whether or not further substituted in the dole ring to any extent and whether or not substituted in the naphthyl ng to any extent, including without limitation the following:	**
	(A)	VH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**
		VH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	
	(C)	VH-018, or 1-Propyl-3-(1-naphthoyl)indole;	
		VH-019, or 1-Hexyl-3-(1-naphthoyl)indole;	
	(E)	VH-073, or 1-Butyl-3-(1-naphthoyl)indole;	**
	(F)	VH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(G)	VH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(H)	VH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl)indole;	**
	(I)	VH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole;	**
	(J)	VH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;	**
	(K)	VH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;	**
	(L)	VH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole;	**
	(M)	AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole;	**
		AM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1-phthalenyl)-methanone;	**
		AM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)-ethanone; and	**
		HJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-)methanone;702	24-(7-2019)
(iii)		hthylmethylindoles, or any compound structurally derived from an H-dol-3-yl-(1-naphthyl) methane by substitution at the nitrogen atom of the dole 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 naphthyl ring to any extent, including without limitation the following:	**
(	(A) JWH-175, or 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane; and	**
(	(B) JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1-naphthyl)methane;	**
(iv)	Naphthoylpyrroles, or 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 and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-307, or (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-ylmethanone;	**
(v)	Naphthylmethylindenes, or any compound structurally derived from 1-(1-napthylmethyl)indene with substitution at the 3-position of the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-176, or E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane;	**
(vi)	Phenylacetylindoles, or 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 and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
(	(A) JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;	**
	(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; and	
	(E) RCS-8, or 1-(2-cyclohexylethyl)-3-(2- methoxyphenylacetyl)indole;	
	Cyclohexylphenols, or 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 without limitation the following:	
(	(A) CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]- phenol;	**

(B) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and	**
(C) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;	**
(viii) Benzoylindoles, or any compound structurally derived from 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 without limitation the following:	**
(A) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;	
(B) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;	**
(C) WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;	**
(D) AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole; and	**
(E) RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1H-indol-3-yl)-methanone;	**
(ix) Adamantoylindoles, or Adamantoylindazoles, including Adamantyl Carboxamide Indoles and Adamantyl Carboxamide Indazoles, or any compound structurally derived from 3-(1-adamantoyl) indole, 3-(1-adamantoyl) indazole, or 3-(2-adamantoyl)indole by substitution at a nitrogen atom of the indole or indazole ring with alkyl, haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl,cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole or indazole ring to any extent and whether or not substituted in the adamantly ring to any extent, including without limitation the following:	**
(A) AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;	**
(B) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;	**
(C) JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide [other name(s): 2NE1, moved in Schedule VI in 2020]	**
(D) AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide;	**
(E) 5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7049**
(F) STS-135 or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide:	**

(x)	Tetramethylcyclopropylcarbonylindoles or any compound structurally derived from 3-(2,2,3,3-tetramethylcyclopropylcarbonyl) indole by substitution at the nitrogen atom of the indole ring with alkyl,haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl, cycloalkylethyl, (N-methylpiperidin-2-yl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole ring to any extent, including without limitation the following:	**
(	(A) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(B) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(C) A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(D) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(E) 5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and	**
(	(F) A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(xi)	Unclassified Synthetic Cannabinoids, including without limitation the following:	**
	(A) CP 50556-1 hydrochloride, 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;	**
(	(B) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;	**
(	(C) 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;	**
(	(D) Dimethylheptylpyran or DMHP;	**
(	(E) WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone;	**
(	(F) URB597, or [3-(3-carbamoylphenyl)phenyl] N-Cyclohexylcarbamate;	**
(	(G) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one;	**
	(H) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]-methanone;	**
	a) URB602, or cyclohexyl N-(3-phenylphenyl)carbamate;	**
(	(I) PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3-carboxylate;	**
(	(J) 5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;	**

(K) BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate; -		**
(L) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide;		**
(M) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide;		**
(N) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L-valine methyl ester	7033-(	9-2018)
(O) MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)-L-valinate	7044-(	(9-2018)
(P) 5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-carboxamido)-3,3-dimethylbutanoate; 7	034-(1	1-2018)
(Q) 5-Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate7	041-(1	1-2018)
(R) MDMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H- indole-3-carboxamido)-3,3-dimethylbutanoate; 7	042-(1	1-2018)
(S) FUB-AMB, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate; 7	021-(1	1-2018)
(T) MDMB-FUBINACA, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate; 7	020-(1	1-2018)
(U) AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-caroboxamide;	7023-(	(7-2019)
(V) AB-CHMINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1- (cyclohexylmethyl)-1H-indazole-3-carboxamide;	7031-(	(7-2019)
(W) MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide;	(1	1-2014)
(X) AB-FUBINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide;	(	(9-2018)
(Y) ADB-PINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide;		(9-2018)
(Z) 5F-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7083-(	(6-2020)
(AA) ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide	7010-(	(4-2021)
(BB) 4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	7043-(	(4-2021)
(CC) 5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7025-(	(5-2022)
(DD) 4-CN-CUMYL-BUTINACA, or 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide 7	'089 –(	(5-2022)

(1	EE) 5F-CUMYL-P7AICA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide	7085-(5-2022)
(I	FF) NM2201, or Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate	7221-(5-2022)
(0	GG) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	7036
(1	HH) FUB-144, or (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone;	7014
(1	I) FUB-AKB48, or N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide;	
(J	J) MDMB-4en-PINACA, or Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate;	
(I	KK) CH-PIATA, or N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide;	
<u>(I</u>	LL) ADB-BUTINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1- butyl-1H-indazole-3-carboxamide	7027
(6) A syı	nthetic substance, derivative, or its isomers with:	
(i)	Similar chemical structure to any substance described in subdivisions (a)(1)-(a)(5) of this section; or	
(ii)	Similar pharmacological effects to any substance described in subdivisions (a)(1)-(a)(5) of this section.	**
	e, except as provided under subsection (c) of this section, the Secretary shall ned substance listed in this section from Schedule VI.	ot delete a

- (b)
- (c) A prescription drug approved by the United States Food and Drug Administration under 21 U.S.C. § 355 is excluded from Schedule VI unless the secretary objects under § 5-64-201.
  - \*-Scheduled before April, 1979.
  - \*\*-Schedule VI is revised to conform to Act 329 of 2013.
  - \*\*\* Schedule VI is revised further to conform to Act 629 of 2023. Each substance added to the Controlled Substances List pursuant to Act 629 of 2023 shall have the following effective dates:
    - (a) For persons who are under twenty-one (21) years of age, the effective date shall be the effective date of Act 629 of 2023; and,
    - (b) For persons who are twenty-one (21) years of age or older, the effective date shall be August 1, 2023.

\*\*\*\* Pursuant to ongoing litigation, and a preliminary injunction against enforcing Act 629 of 2023, the changes made to the List of Controlled Substances pursuant to Act 629 of 2023 are not enforceable until a final order issued in the matter, Bio Gen, LLC, et al. v. Sarah Huckabee Sanders, et al., Case No. 4:23-CV-00718-BRW, Central Division, Eastern District of Arkansas, United States District Court.

# QUESTIONNAIRE FOR FILING PROPOSED RULES WITH THE ARKANSAS LEGISLATIVE COUNCIL

	ARTMENT		
	ARD/COMMISSION		
BOA	RD/COMMISSION DIRECTOR		
CON	VTACT PERSON		
ADD	ORESS		
	PHONE NO EMAIL		
NAN	ME OF PRESENTER(S) AT SUBCOMMITTEE MEETING		
PRE	SENTER EMAIL(S)		
	<u>INSTRUCTIONS</u>		
Ques what	rder to file a proposed rule for legislative review and approval, please submit this Legislative stionnaire and Financial Impact Statement, and attach (1) a summary of the rule, describing the rule does, the rule changes being proposed, and the reason for those changes; (2) both a kup and clean copy of the rule; and (3) all documents required by the Questionnaire.		
of Re	If the rule is being filed for permanent promulgation, please email these items to the attention of Rebecca Miller-Rice, <u>miller-ricer@blr.arkansas.gov</u> , for submission to the Administrative Rules Subcommittee.		
Dire	e rule is being filed for emergency promulgation, please email these items to the attention of ctor Marty Garrity, <a href="mailto:garritym@blr.arkansas.gov">garritym@blr.arkansas.gov</a> , for submission to the Executive committee.		
Pleas	se answer each question completely using layman terms.		
**** 1.	**************************************		
2.	What is the subject of the proposed rule?		
3.	Is this rule being filed under the emergency provisions of the Arkansas Administrative Procedure Act? Yes No		
	If yes, please attach the statement required by Ark. Code Ann. § 25-15-204(c)(1).		
	If yes, will this emergency rule be promulgated under the permanent provisions of the Arkansas Administrative Procedure Act? Yes No		

4.	Is this rule being filed for permanent promulgation? Yes No
	If yes, was this rule previously reviewed and approved under the emergency provisions of the Arkansas Administrative Procedure Act? Yes No
	If yes, what was the effective date of the emergency rule?
	On what date does the emergency rule expire?
5.	Is this rule required to comply with a <i>federal</i> statute, rule, or regulation? Yes No
	If yes, please provide the federal statute, rule, and/or regulation citation.
_	
6.	Is this rule required to comply with a <i>state</i> statute or rule? Yes No
	If yes, please provide the state statute and/or rule citation.
7.	Are two (2) rules being repealed in accord with Executive Order 23-02? Yes No
	If yes, please list the rules being repealed.
	If no, please explain.
8.	Is this a new rule? Yes No
	Does this repeal an existing rule? Yes No If yes, the proposed repeal should be designated by strikethrough. If it is being replaced with a new rule, please attach both the proposed rule to be repealed and the replacement rule.
	Is this an amendment to an existing rule? Yes No If yes, all changes should be indicated by strikethrough and underline. In addition, please be

sure to label the markup copy clearly as the markup.

9.	What is the state law that grants the agency its rulemaking authority for the proposed rule, outside of the Arkansas Administrative Procedure Act? Please provide the specific Arkansas Code citation(s), including subsection(s).
10.	Is the proposed rule the result of any recent legislation by the Arkansas General Assembly? Yes No
	If yes, please provide the year of the act(s) and act number(s).
11.	What is the reason for this proposed rule? Why is it necessary?

Please provide the web address by which the proposed rule can be accessed by the public as provided in Ark. Code Ann. § 25-19-108(b)(1).
Will a public hearing be held on this proposed rule? Yes No
If yes, please complete the following:
Date:
Time:
Place:
e be sure to advise Bureau Staff if this information changes for any reason.
On what date does the public comment period expire for the permanent promulgation of the rule? Please provide the specific date.
What is the proposed effective date for this rule?
Please attach (1) a copy of the notice required under Ark. Code Ann. § 25-15-204(a)(1) and (2) proof of the publication of that notice.
Please attach proof of filing the rule with the Secretary of State, as required by Ark. Code Ann. § 25-15-204(e)(1)(A).
Please give the names of persons, groups, or organizations that you anticipate will comment on these rules. Please also provide their position (for or against), if known.
Is the rule expected to be controversial? Yes No  If yes, please explain.

# FINANCIAL IMPACT STATEMENT

# PLEASE ANSWER ALL QUESTIONS COMPLETELY.

DEP	PARTMENT	
	ARD/COMMISSION	
PER	SON COMPLETING THIS STATEMENT	
TEL	EPHONE NO. EMAIL	
To comply with Ark. Code Ann. § 25-15-204(e), please complete the Financial Impact Statement and email it with the questionnaire, summary, markup and clean copy of the rule, and other documents. Please attach additional pages, if necessary.		
TITI	LE OF THIS RULE	
1.	Does this proposed, amended, or repealed rule have a financial impact? Yes No	
2.	Is the rule based on the best reasonably obtainable scientific, technical, economic, or other evidence and information available concerning the need for, consequences of, and alternatives to the rule?  Yes  No	
3.	In consideration of the alternatives to this rule, was this rule determined by the agency to be the least costly rule considered? Yes No	
	If no, please explain:	
	(a) how the additional benefits of the more costly rule justify its additional cost;	
	(b) the reason for adoption of the more costly rule;	
	(c) whether the reason for adoption of the more costly rule is based on the interests of public health, safety, or welfare, and if so, how; and	
	(d) whether the reason for adoption of the more costly rule is within the scope of the agency's statutory authority, and if so, how.	
4.	If the purpose of this rule is to implement a <i>federal</i> rule or regulation, please state the following	

(a) What is the cost to implement the federal rule or regulation?

	Next Fiscal Year
General Revenue	General Revenue
Federal Funds	Federal Funds
Cash Funds	Cash Funds
Special Revenue	Special Revenue_
Other (Identify)	Other (Identify)
Total	Total
(b) What is the additional cost of the	state rule?
Current Fiscal Year	Next Fiscal Year
General Revenue	General Revenue
Federal Funds	Federal Funds
Cash Funds	Cash Funds
Special Revenue	Special Revenue_
Other (Identify)	Other (Identify)
Total	Total
business subject to the proposed, ame rule, and explain how they are affecte <u>Current Fiscal Year</u>	ended, or repealed rule? Please identify those subject
business subject to the proposed, ame rule, and explain how they are affecte Current Fiscal Year  \$	ended, or repealed rule? Please identify those subjected.  Next Fiscal Year  \$
business subject to the proposed, ame rule, and explain how they are affecte Current Fiscal Year  \$  What is the total estimated cost by fiscal in the state of the proposed, ame rule, and explain how they are affected and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed, ame rule, and explain how they are affected in the proposed in the propos	

7. With respect to the agency's answers to Questions #5 and #6 above, is there a new or increased cost or obligation of at least one hundred thousand dollars (\$100,000) per year to a private individual, private entity, private business, state government, county government, municipal government, or to two (2) or more of those entities combined?

Yes No

If yes, the agency is required by Ark. Code Ann. § 25-15-204(e)(4) to file written findings at the time of filing the financial impact statement. The written findings shall be filed simultaneously with the financial impact statement and shall include, without limitation, the following:

- (1) a statement of the rule's basis and purpose;
- (2) the problem the agency seeks to address with the proposed rule, including a statement of whether a rule is required by statute;
- (3) a description of the factual evidence that:
  - (a) justifies the agency's need for the proposed rule; and
  - (b) describes how the benefits of the rule meet the relevant statutory objectives and justify the rule's costs:
- (4) a list of less costly alternatives to the proposed rule and the reasons why the alternatives do not adequately address the problem to be solved by the proposed rule;
- (5) a list of alternatives to the proposed rule that were suggested as a result of public comment and the reasons why the alternatives do not adequately address the problem to be solved by the proposed rule;
- (6) a statement of whether existing rules have created or contributed to the problem the agency seeks to address with the proposed rule and, if existing rules have created or contributed to the problem, an explanation of why amendment or repeal of the rule creating or contributing to the problem is not a sufficient response; and
- (7) an agency plan for review of the rule no less than every ten (10) years to determine whether, based upon the evidence, there remains a need for the rule including, without limitation, whether:
  - (a) the rule is achieving the statutory objectives;
  - (b) the benefits of the rule continue to justify its costs; and
  - (c) the rule can be amended or repealed to reduce costs while continuing to achieve the statutory objectives.



# SUMMARY OF PROPOSED AMENDMENTS TO RULES PERTAINING TO THE LIST OF CONTROLLED SUBSTANCES FOR THE STATE OF ARKANSAS

The proposed listed amendments update List of Controlled Substances to include these drugs.

**GOVERNOR** 

- 1) Meta-fluorofentanyl [other name(s): N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4yl)propionamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (RR).
- 2) Meta-fluoroisobutyryl fentanyl [other name(s): N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4yl)isobutyramide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (SS).
- 3) Para-methoxyfuranyl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (TT).
- 4) Para-methylcyclopropyl fentanyl [other name(s): N-(4-methylphenyl)-N-(1phenethylpiperidin-4-yl)cyclopropanecarboxamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (UU).
- 5) 3-furanyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3carboxamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (VV).
- 6) 2',5'-dimethoxyfentanyl [other name(s): N-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-Nphenylpropionamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (WW).
- 7) Isovaleryl fentanyl [other name(s): 3-methyl-N-(1-phenethylpiperidin-4-yl)-Nphenylbutanamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 5, Schedule I, (b), (51), (vi), (XX).



- 8) Ortho-fluorofuranyl fentanyl [other name(s): *N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 6, Schedule I, (b), (51), (vi), (YY).
- 9) Alpha'-methyl butyryl fentanyl [other name(s): 2-methyl-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbutanamide]. The DEA has placed this opioid analgesic into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 6, Schedule I, (b), (51), (vi), (ZZ).
- 10) Etodesnitazene [other name(s): Etazene], *N*-Pyrrolidino Etonitazene (other name(s) Etonitazepyne), and Protonitazene are Schedule I controlled substances. To follow DEA, controlled substance code numbers have been set forth opposite of these substances. Page 6, Schedule I, (b), (54), (vi), (E), Page 6, Schedule I, (b), (54), (vi), (O), and Page 6, Schedule I, (b), (54), (vi), (Q).
- 11) 2-Methyl AP–237. [other name(s): 1-(2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl)butan-1-one]. The DEA has placed this synthetic opioid into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Page 7, Schedule I, (b), (55).
- 12) 3-methylmethcathinone (other names: 3–MMC). The DEA has identified this synthetic cathinone as a positional isomer of mephedrone a currently listed controlled substance. This drug without a recognized medical use would be included as Schedule I with subsequent numbering changes to follow. Page 13, Schedule I, (f), (2), (xvii).
- 13) 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one [other names: *alpha*-PiHP, and Alpha-Pyrrolidinoisohexanophenone). The DEA has identified this synthetic cathinone as a positional isomer of Alpha-PHP a currently listed controlled substance. This drug without a recognized medical use would be included as Schedule I. Page 14, Schedule I, (f), (2), (xxviii).
- 14) The following items are marked for clean up:
  - a. Page 2, Schedule I, (b), (46);
  - b. Page 3, Schedule I, (b), (51), (vi), (B);
  - c. Page 3, Schedule I, (b), (51), (vi), (C);
  - d. Page 3, Schedule I, (b), (51), (vi), (G);
  - e. Page 3, Schedule I, (b), (51), (vi), (I);
  - f. Page 4, Schedule I, (b), (51), (vi), (N);
  - g. Page 12, Schedule I, (f), (1), (iv);
  - h. Page 12, Schedule I, (f), (1), (ix);
  - i. Page 12, Schedule I, (f), (1), (xi);
  - j. Page 13, Schedule I, (f), (2), (iii);
  - k. Page 17. Schedule II, (d), (3);
  - 1. Page 19, Schedule III, (c), (13);



- m. Page 23, Schedule III, (g), (51); and
- n. Page 27, Schedule IV, (d), (1).
- 15) Xylazine The potential adverse health effects when abused, and increasing national prevalence of xylazine utilized as an adulterating agent to other illicit substances poses a threat to public health and safety. The substance will be included as a Schedule III controlled substance, Page 19, Schedule III, (c), (15), (i through vi), with outlined exceptions utilizing the following language:

Xylazine and any material, compound, mixture, or preparation which contains any quantity of xylazine, 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, except in the following uses:

- i) Dispensing, prescribing, or administering, to an animal, a drug containing xylazine that has been approved by the United States Secretary of Health and Human Services under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b);
- ii) Dispensing, prescribing, or administering xylazine to an animal that is permissible under section 512 (a)(4) of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(a)(4));
- iii) Possessing a drug containing xylazine, as described in this Section (15), for animal use:
  - (A). By a licensed pharmacist or licensed veterinarian; or
  - (B). Pursuant to a valid prescription from a licensed veterinarian.
- iv) Possessing, manufacturing, distributing, or using xylazine as an active pharmaceutical ingredient for manufacturing an animal drug either:
  - (A). Approved under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b); or
  - (B). Issued an investigation use exemption under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(j));
- Manufacturing, distributing, or using a xylazine bulk chemical for pharmaceutical compounding by a licensed pharmacist or veterinarian; or
- vi) Another use approved or permissible under the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 301, et seq.) or under 21 CFR Part 530, Subpart B.
- 16) Zuranolone. The FDA approved this drug for use in treatment of post-partum depression. This drug would be included as Schedule IV to follow DEA. Page 27, Schedule IV, (c), (61).



17) ADB–BUTINACA or *N*-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1*H*-indazole-3-carboxamide. The DEA has identified this synthetic cannabinoid as positional isomer of AB-PINACA, a currently listed controlled substance. This drug without a recognized medical use would be included as Schedule VI. Page 36, Schedule VI, (a), (5), (xi), (LL).

List Of Controlled Substances



For the State Of Arkansas Pursuant to the provisions of Arkansas Code Annotated § 5-64-201 and § 5-64-216 of the laws of the State of Arkansas, the Secretary of the Arkansas Department of Health or duly authorized agent, as specified by law, is giving public notice of the publication of the List of Controlled Substances for the State of Arkansas.

Due consideration has been given applicable federal regulations, current scientific knowledge regarding the listed substances, the evidence of actual or relative potential for abuse, the history and current patterns of abuse, the risk to the public health, and potential to produce psychic or psychological dependence liability.

Based on these considerations the attached listing of the Schedule of Controlled Substances and the corresponding drugs that are included in each schedule is hereby promulgated by the Secretary of the Arkansas Department of Health as the List of Controlled Substances for the State of Arkansas.

Each controlled substance or basic class thereof has been assigned an "Administration Controlled Substance Code Number" for purposes of identification. These numbers are for internal management and are used as a means to identify substances with complex and cumbersome chemical names.

Next to the code number is the date the substance was placed in schedule by the Secretary of the Arkansas Department of Health. A "\*" denotes the substance was scheduled prior to April, 1979.

I, Shane David, Pharm.D., Branch Chief, Health Systems Licensing and Certifications for the Arkansas Department of Health, do hereby certify that the documents attached hereto are true and correct copies of the current List of Controlled Substances adopted by the Arkansas State Board of Health in accordance with Arkansas state law.

	Shane David, Pharm.D.
STATE OF ARKANSAS ) COUNTY OF SALINE )	
I, Marci Middleton-Yates, do he before me and signed the above Sworn and subscribed to	
	Notary Public
My commission expires	

#### ARKANSAS DEPARTMENT OF HEALTH

## LIST OF CONTROLLED SUBSTANCES

#### **SECTION I AUTHORITY**

The following scheduling of these controlled substances has been hereby promulgated pursuant to Arkansas Code Annotated §5-64-201 and §5-64-216.

#### **SECTION II PURPOSE**

Due consideration has been given applicable Federal regulations, current scientific knowledge regarding the listed substances, the evidence of actual or relative potential for abuse, the history and current patterns of abuse, the risk to the public health, and potential to produce psychic or psychological dependence liability.

## SECTION III GENERAL REQUIREMENTS

(Attached copy of the listing of scheduling of controlled substances)

## **SECTION IV REPEAL**

All lists of schedules of controlled substances in conflict herewith are hereby repealed.

#### CERTIFICATION

This will certify the following list of scheduling of controlled substances was adopted by the Arkansas State Board of Health at a session of the Board held in Little Rock, Arkansas on the \_\_\_\_ day of October, 2024, and after a Public Hearing on the, held in Little Rock, Arkansas, at the State Department of Health Building.

Jennifer Dillaha, MD Secretary of Arkansas State Board of Health Director of the Arkansas Department of Health

#### **ARTICLE II**

# **SCHEDULE I**

- (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 DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Opiates: (Narcotic Drugs)</u> 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, salts is possible within the specific chemical designation (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

(1) A	cetylmethadol	9601*
(2) A	llylprodine	9602*
(3) A	lphacetylmethadol (except Levo-alphacetylmethadol (LAAM)	-9603*
(4) A	lphameprodine	9604*
(5) A	lphamethadol	9605*
(6) B	enzethidine	9606*
(7) B	etacetylmethadol	9607*
(8) B	etameprodine	9608*
(9) B	etamethadol	9609*
(10)	Betaprodine	9611*
(11)	Dextromoramide	9613*
(12)	Diampromide	9615*
(13)	Diethylthiambutene	9616*
(14)	Difenoxin	9168*
(15)	Dimenoxadol	9617*
(16)	Dimepheptanol	9618*
(17)	Dimethylthiambutene	9619*
(18)	Dioxaphetyl butyrate	9621*
(19)	Dipipanone	9622*
(20)	Ethylmethylthiambutene	9623*
(21)	Etoxeridine	9625*
(22)	Furethidine	9626*
(23)	Hydroxypethidine	9627*

(24)	Ketobemidone	9628*
(25)	Levomoramide	9629*
(26)	Levophenacylmorphan	9631*
(27)	Morpheridine	9632*
(28)	MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)]	0661-(10-1985)
(29)	Noracymethadol	9633*
(30)	Norlevorphanol	9634*
(31)	Normethadone	9635*
(32)	Norpipanone	9636*
(33)	PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] - 9	0663-(10-1985)
(34)	Phenadoxone	9637*
(35)	Phenampromide	9638*
(36)	Phenomorphan	9647*
(37)	Phenoperidine	9641*
(38)	Piritramide	9642*
(39)	Proheptazine	9643*
(40)	Properidine	9644*
(41)	Propiram	9649*
(42)	Racemoramide	9645*
(43)	Tilidine	9750-(9-1981)
(44)	Trimeperidine	9646*
(45)	Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide]	(4-2017)
(46) d	AH-7921 [other name(s): 3,4-dichloro-N-[(1-limethylamino)cyclohexylmethyl]benzamide]	9551-(4-2017)
	W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4- chlorophenyl)sulfonamide]	(4-2017)
	W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4- chlorophenyl)Sulfonamide]	(4-2017)
(49)	MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine]	9560-(4-2017)
(50) N	U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)-N-methylbenzamide]	9547-(4-2017)
	Fentanyl-related substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Fentanyl-related substance means any substance not otherwise listed, and for which no exemption or approval is in effect under section	

	of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that is turally related to fentanyl by one or more of the following modifications:	(4-2021)
(i)	Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;	
(ii)	Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iii)	Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iv)	Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; or	
(v)	Replacement of the N-propionyl group by another acyl group.	
(vi)	Fentanyl-related substances shall include, but are not limited to:	
(4	A) Acetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	9815-(2-1986)
(1	B) Alpha-methylfentanyl [other name(s): (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propronanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine)]	- 9814-(6-1982)
(0	C) Alpha-methylthiofentanyl [other name(s): N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide]	
(1	D) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	- 9830-(2-1986)
(1	E) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide]	- 9831-(2-1986)
(1	F) 3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-piperidyl]-N-Phenylpropanamide)]	9813-(10-1985)
(0	G)3-methylthiofentanyl [other name(s): N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide]	- 9833-(2-1986)
(1	H) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide]	9812-(11-1986)
(]	Thiofentanyl [other name(s): (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide]	9835-(2-1986)
(.	(1) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9821-(4-2017)
(1	K) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	9822-(4-2017)
(1	L) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2 yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	-

(M) Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2 (4-methylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	
(N) Valeryl fentanyl [other name(s): N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide]	9840-(4-2017)
(O) Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran-2-carboxamide]	
(P) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	9827-(4-2017)
(Q) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	9838-(4-2017)
(R) 4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	
(S) Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide]	
(T) Acryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]	9811-(6-2020)
(U)4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9824-(6-2020)
(V) Tetrahydrofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-y N-phenyltetrahydrofuran-2-carboxamide]	
(W) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-y N-phenylcyclopropanecarboxamide]	l)- 9845-(4-2021)
(X) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9825-(4-2021)
(Y) Ortho-fluorofentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9816-(4-2021)
(Z) Crotonyl fentanyl [other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide]	9844-(5-2022)
(AA) Cyclopentyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl N-phenylcyclopentanecarboxamide]	
(BB) Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl) N-(1-phenethylpiperidin-4-yl)isobutyramide]	9826-(5-2022)
(CC) Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	9837-(5-2022)
(DD) Beta-methyl fentanyl [other name(s): N-phenyl-N-(1-(2-phenylpropyl) piperidin-4-yl)propionamide]	
(EE) Beta'-phenyl fentanyl [other name: N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide]	9842-(5-2022)

· /	2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2-pphenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	9855-(5-2022)
	4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) din-4-yl)-N-phenylacetamide]	9819-(5-2022)
	Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-enethylpiperidin-4-yl)butyramide]	9846-(5-2022)
(II) Ortho phene	-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-ethylpiperidin-4-yl)acetamide]	-9848-(5-2022)
	-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2-ylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9820-(5-2022)
	Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1-ethylpiperidin-4-yl)propionamide]	9817-(5-2022)
	Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-vlbenzamide]	9841-(5-2022)
	Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-enylthiophene-2-carboxamide]	
(NN) yl)(ph	Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4-nenyl)carbamate]	9851-(5-2022)
(OO) phene	Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-ethylpiperidin-4-yl)acrylamide]	9852-(5-2022)
(PP) N-(1-)	Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-phenethylpiperidin-4-yl)isobutyramide]	9853-(5-2022)
	Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-enethylpiperidin-4-yl)furan-2-carboxamide]	9854-(5-2022)
(RR) phene	Meta-fluorofentanyl [other name(s): N-(3-fluorophenyl)-N-(1-ethylpiperidin-4-yl)propionamide]	9857
	Meta-fluoroisobutyryl fentanyl [other name(s): N-(3-fluorophenyl)-phenethylpiperidin-4-yl)isobutyramide	9858
	Para-methoxyfuranyl fentanyl [other name(s): N-(4-oxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	9859
	Para-methylcyclopropyl fentanyl [other name(s): N-(4-ylphenyl)-N-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide]	9865
	3-furanyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-ylfuran-3-carboxamide]	9860
` /	2',5'-dimethoxyfentanyl [other name(s): N-(1-(2,5-hoxyphenethyl)piperidin-4-yl)-N-phenylpropionamide]	9861
` /	Isovaleryl fentanyl [other name(s): 3-methyl-N-(1-ethylpiperidin-4-yl)-N-phenylbutanamide]	9862

	(YY) Ortho-fluorofuranyl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	9863
	(ZZ) Alpha'-methyl butyryl fentanyl [other name(s): 2-methyl-N-(1-	
	phenethylpiperidin-4-yl)-N-phenylbutanamide]	
(52)	• •	
(53)	Brorphine	9098
S	Benzimidazole-opioid substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Benzimidazole-opioid substances includes any substance, not otherwise listed or excepted, 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 structurally has benzimidazole ring with an ethylamine at its 1-position and benzyl group at its 2-position:	
(	(i) With or without substitution on the benzimidazole;	
(	(ii) With or without substitution at the ethylamine;	
(	(iii) With or without inclusion of the ethylamine in a cyclic structure;	
(	(iv) With or without substitution on the benzyl ring; or	
(	(v) With or without replacement of the benzyl ring with an aromatic ring.	
(	(vi) Benzimidazole-opioid substances shall include but are not limited to:	
	(A)4'-Hydroxy Nitazene;	
	(B) 5-Aminoisotonitazene;	
	(C) Butonitazene;	
	(D)Clonitazene;	9612 *
	(E) Etodesnitazene, [other name(s): Etazene];	9765
	(F) Etonitazene;	9624*
	(G) Flunitazene;	
	(H) Isotonitazene;	9614
	(I) Isotodesnitazene;	
	(J) Metodesnitazene;	
	(K) Metonitazene;	9757
	(L) N-Desethyl Etonitazene;	
	(M) N-Desethyl Isotonitazene;	
	(N) N-Piperidinyl Etonitazene [other name(s): Etonitazepipne];	
	(O) N-Pyrrolidino Etonitazene [other name(s): Etonitazepyne];	9758
	(P) N-Pyrrolidino Protonitazene; and	
	(Q) Protonitazene	9759

		2-Methyl AP–237 [other name(s): 1-(2-methyl-4-(3-phenylprop-2-en-1-l)piperazin-1-yl)butan-1-one]	9664
(c)	sched the ex	m derivatives: (Narcotic Drugs) Unless specifically excepted or unless listed in ano lule, any of the following opium derivatives, its salts, isomers, and salts of isomers what stence of such salts, isomers, and salts of isomers is possible within the specific chernation:	nenever
	(1) A	cetorphine	9319*
	(2) A	cetyldihydrocodeine	9051*
	(3) B	enzylmorphine	9052*
	(4) C	odeine methylbromide	9070*
	(5) C	odeine-N-Oxide	9053*
	(6) C	yprenorphine	9054*
	(7) D	esomorphine	9055*
	(8) D	ihydromorphine	9145*
	(9) D	rotebanol	9335*
	(10)	Etorphine (except hydrochloride salt)	9056*
	(11)	Heroin	9200*
	(12)	Hydromorphinol	9301*
	(13)	Methyldesorphine	9302*
	(14)	Methyldihydromorphine	9304*
	(15)	Morphine methylbromide	9305*
	(16)	Morphine methylsulfonate	9306*
	(17)	Morphine-N-Oxide	9307*
	(18)	Myrophine	9308*
	(19)	Nicocodeine	9309*
	(20)	Nicomorphine	9312*
	(21)	Normorphine	9313*
	(22)	Pholcodine	9314*
	(23)	Thebacon	9315*
	(24)	Mitragynine	(11-2015)
	(25)	7-Hydroxymitragynine	(11-2015)

(d) **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 substance, 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 this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

(1) alpha-ethyltryptamine	7249-(12-1993)
Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-i ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET.	ndole-3-
(2) 4-bromo-2,5-dimethoxy-amphetamine	7391*
Some trade or other names: 4-bromo-2,5-dimethoxy-alphamethylphenethylamine; 4-bromo-2,5-DMA.	
(3) 4-bromo-2,5-dimethoxyphenethylamine	7392-(8-1995)
Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1 an alpha-desmethyl DOB; 2C-B, Nexus.	ninoethane;
(4) 2,5-dimethoxyamphetamine	7396*
Some trade or other names: 2,5-dimethoxy-alpha-methylphenethyla DMA.	mine; 2,5-
(5) 2,5-dimethoxy-4-ethylamphetamine	7399-(3-1988)
Some trade or other names: DOET.	
(6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine	7348-(1-2005)
Some trade or other names: 2C-T-7.	
(7) 4-methoxyamphetamine	7411*
Some trade or other names: 4-methoxy-alpha- methylphenethylami paramethoxyamphetamine; PMA.	ne;
(8) 5-methoxy-3,4-methylenedioxy-amphetamine	7401*
(9) 4-methyl-2,5-dimethoxyamphetamine	7395*
Some trade and other names: 4-methyl-2,5-dimethoxy-alphamethylphenethylamine; "DOM"; and "STP".	
(10) 3,4-methylenedioxy amphetamine	7400*
(11) 3,4-methylenedioxymethamphetamine	7405-(10-1985)
Some trade or other names: MDMA)	
(12) 3,4-methylenedioxy-N-ethylamphetamine	7404-(6-1990)
Some trade or other names: N- ethy-alpha-methyl-3,4 (methylenedic phenethylamine, N-ethyl MDA; MDE; MDEA.	oxy)
(13) N-hydroxy-3,4-methylenedioxyamphetamine	7402-(6-1990)
Some trade or other names: N-hydroxy-alpha-methyl-3,4(methylene phenethylamine; N-hydroxy MDA	edioxy)

(14)	3,4,5-trimethoxy amphetamine	7390*
(15)	5-methoxy-n,n-dimethyltryptamine 5-MeO-DMT	- 7431*(1-2011)
(16)	alpha-methyltryptamine	- 7432-(7-2005)
	Some trade or other names: AMT	
(17)	Bufotenine	7433*
	Some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine.	
(18)	Diethyltryptamine	7434*
	Some trade or other names: N,N-Diethyltryptamine;DET.	
(19)	Dimethyltryptamine	7435*
	Some trade or other names: DMT	
(20)	5-methoxy-N,N-diisopropyltryptamine	- 7439-(7-2005)
	Some trade or other names: 5-MeO-DIPT.	
(21)	Ibogaine	7260*
	Some trade and other names: 7-Ethyl-6,6 beta; 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.	
(22)	Lysergic acid diethylamide	7315*
(23)	Mescaline	7381*
(24)	Parahexyl	- 7374-(7-1983)
	Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl.	
(25)	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, salts, derivative, mixture or preparation of such plant, its seeds or extracts. (Interprets 21 USC 812 (c), Schedule I (c) (12)).	
(26)	N-ethyl-3-piperidyl benzilate	7482*
(27)	N-methyl-3-piperidyl benzilate	7484*
(28)	Psilocybin	7437*
(29)	Psilocyn	7438*
(30)	Ethylamine Analog of phencyclidine	7455*

	Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE.	
(31)	Pyrrolidine Analog of phencyclidine	7458*
	Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCPy; PHP	
(32)	Thiophene Analog of phencyclidine	7470*
	Some trade or other names: 1-[1-(2-thienyl) cyclohexyl] Piperidine; 2-Thienyl analog of phencyclidine; TPCP; TCP.	
(33)	1-[1-(2-Thienyl)cylcohexyl]pyrrolidine	- 7473-(9-1989)
	Some other trade or other names: TCPy.	
(34)	N,N-Diallyl-5-Methoxytryptamine;	(6-2012)
	Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT	
(35)	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 5C-NBOMe	
	2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5I-NBOMe	(8-2013)
(37)	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E	7509-(11-2013)
(38)	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D	7508-(11-2013)
(39)	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C	7519-(11-2013)
(40)	2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I	7518-(11-2013)
(41)	2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-2	7385-(11-2013)
(42)	2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-4	7532-(11-2013)
(43)	2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H	7517-(11-2013)
(44)	2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N	7521-(11-2013)
(45)	2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine; 2C-P	7524-(11-2013)
	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine; 25B-	(9-2018)
	2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25B-	(11-2018)
(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH	(11-2018)
	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5E-NBOMe	(7-2019)
	2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 25H-NBOMe	(7-2019)
(51)	2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]- phenol; 25C-	

	(52)	2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25H-NBOH	(7-2019)
	(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	1245-(5-2022)
		Some trade or other names: Para-methoxymethamphetamine; PMMA	
	(54)	2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one	7286
		Some trade or other names: Methoxetamine; MXE	
(e)	com depr	pressants: Unless specifically excepted or unless listed in another schedule, any napound, mixture, or preparation which contains any quantity of the following substressant effect on the central nervous system, including its salts, isomers, and salts enever the existence of such salts, isomers, and salts of isomers is possible within the mical designation:	tances having a of isomers
	i	gamma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate], and its known precursors and analogs. Precursors include but are not limited to: gamma-butyrolactone	
	(2)	Mecloqualone	2572*
	(3)		2565*
	1 1 1	Benzodiazepine substances, their isomers, esters, ethers, salts and salts of isomers esters and ethers. Benzodiazepine substances includes any substance, not otherwise listed or excepted, 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 structurally has a fused 1,4-diazepine and benzene ring structure with a phenyl connected to the diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or any combination thereof. Benzodiazepine substances shall include but are not limited to:	
	(	(i) Bromazolam;	
	(	(ii) Clonazolam;	(4-2021)
	(	(iii) Flualprazolam;	(4-2021)
	(	(iv) Flubromazepam;	
	(	(v) Flubromazolam;	(4-2021)
	(	(vi) Phenazepam;	(6-2012)
	(	(vii) Phenazolam [other name(s): Clobromazolam].	
	i 1	Thienodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Thienodiazepine substances includes any substance, not otherwise listed or excepted, 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 structurally has a fused 1,4-diazepine and thiophene ring structure with a phenyl connected to the 1,4-diazepine ring, with any substitution(s) or	

		phen	cement(s) on the 1,4-diazepine or thiophene ring, any substitution(s) on the yl ring, or any combination thereof. Thienodiazepine substances shall include re not limited to:	
		(i)	Etizolam	(12-2014)
(f)	Sti	mula	nts:	
	(1)	comp subst	ss specifically excepted or unless listed in another schedule, any material, bound, mixture, or preparation which contains any quantity of the following ances having a stimulant effect on the central nervous system, including its isomers, and salts of isomers:	
		(i)	Cathinone	1235-(3-1988)
		(ii)	(±) CIS-4-Methylaminorex [(±)CIS-4,5-dihydro-4- methyl-5-phenyl-2-	4.500 (6.4000)
			oxazolamine]	,
		` /	Fenethylline	` ′
		(iv)	N-Benzylpiperazine [other name(s): BZP, 1-Benzylpiperazine]	7493-(1-2005)
		(v)	N-ethylamphetamine	1475-(6-1982)
		(vi)	N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers	9818-(2-1986)
		(vii)	N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropan-mide (thenylfentanyl), its optical isomers, salts and salts of isomers	9834-(2-1986)
		(viii)	N,N-Dimethylamphetamine [other name(s): N,N,Alpha-trimethylbenzeneethanamine; N,N,Alpha-trimethylphenethylamine], its salts, optical isomers, and salts of optical isomers	1480-(2-1989)
		(ix)	Methcathinone [other name(s):2-Methylamine-Proprophenone, alpha (methylamino)- Proprophenone, 2 (methylamino)-1-phenylpropan-1-one, alpha-N-Methylaminopropiophenone, monomethylpropion, ephedrone, N-methylcathinone, methycathinone, AL-464, AL-422, AL-463 and UR-1431], its salts, optical isomers and salts of optical isomers	237-(12-1993)
		(x)	Aminorex [other name(s): aminoraphen, 2-amino-5 phenyl-2-oxazoline, or 4,5 dihyrdo-5-phenyl-2-oxazolamine], its salts, optical isomers, and salts of optical isomers	.585-(12-1993)
		(xi)	4,4'-Dimethylaminorex [other name(s): 4,4'-DMAR, 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine, or 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine]	1595-(5-2022)
		(xii)	Amineptine	1219
		(xiii)	Mesocarb	1227
		(xiv)	Methyl-N-ethylcathinone (MEC)	(6-2014)

	(XV)	Methiopropamine. [other name(s): N-methyl-1-(thiophen-2-yl)propan-2-amine]	1478
(2)	indirection contact followexiste specific specific contact for the contact followexiste specific specific contact for the cont	material, compound, mixture, or preparation, whether produced directly or ectly from a substance of vegetable origin or independently by means of nical synthesis or by a combination of extraction and chemical synthesis, that any quantity of the following substances, or that contains any of the wing substances' analogs, salts, isomers, and salts of isomers when the ence of the analogs, salts, isomers, and salts of isomers is possible within the fic chemical designation, with the following chemical structure is included it dule I:	
	(i)	4-Methylmethcathinone (Mephedrone)	- 1248-(3-2011)
	(ii)	Methylenedioxypyrovalerone (MDPV)	(3-2011)
	(iii)	3,4-Methylenedioxy-N-methylcathinone (Methylone)	- 7540-(3-2011)
	(iv)	4-Methoxymethcathinone	(3-2011)
	(v)	3-Fluoromethcathinone	(3-2011)
	(vi)	4-Fluoromethcathinone	(3-2011)
	(vii)	1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)	7541-(11-2014)
	(viii)	Alpha-Pyrrolidinopentiophenone (Alpha-PVP)	7545-(11-2015)
	(ix)	4-methyl-N-ethylcathinone (4-MEC)	1249-(9-2018)
	(x)	4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)	7498-(9-2018)
	(xi)	2-(methylamino)-1-phenylpentan-1-one (Pentedrone)	1246-(9-2018)
	(xii)	1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)	7542-(9-2018)
	(xiii)	4-fluoro-N-methylcathinone (4-FMC, Flephedrone)	1238-(9-2018)
	(xiv)	3-fluoro-N-methylcathinone (3-FMC)	1233-(9-2018)
	(xv)	1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)	1258-(9-2018)
	(xvi)	Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)	7546-(9-2018)
	(xvii	3-methylmethcathinone (3–MMC)	1259
	(xvii	i) A compound, unless listed in another schedule or a legend drug, that is structurally derived from 2-Amino-phenyl-1-propanone by modification or by substitution:	
	(.	A) In the phenyl ring to any extent with alkyl, alkoxy, alkylenedioxy, haloalkyl or halide substituents, whether or not further substituted in the phenyl ring by one (1) or more other univalent substituents;	
	(1	B) At the 3-position with an alkyl substituent; or	
	(	C) At the nitrogen atom with alkyl or dialkyl groups, or by inclusion of the nitrogen atom in a cyclic structure.	

	(xix)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one (N-Ethylpentylone)	-7543-(7-2019)				
	(xx)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)	-7547-(4-2021)				
	(xxi)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone)	-7549-(4-2021)				
	(xxi	(i)2-(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone; Alpha-Ethylaminohexanophenone)]	7246				
	(xxi	ii) 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (Alpha-Pyrrolidinohexanophenone; Alpha-PHP)]	7544				
	(xxi	v) 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one [other name(s): (4-Methyl-alpha-ethylaminopentiophenone; 4-MEAP)]	7245				
	(xxv	1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (4'-Methyl-alpha-pyrrolidinohexiophenone; MPHP)]	7446				
	(xxv	i) 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha-Pyrrolidinoheptaphenone; PV8)]	7548				
	(xxv	ii) 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): (4'-Chloro-alpha-pyrrolidinovalerophenone; 4-chloro-alpha-PVP)]	7443				
	(xxv	iii) 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): alpha-PiHP, Alpha-Pyrrolidinoisohexanophenone]					
<u>SC</u>	<u>HEDUL</u>	<u>E II</u>					
(a)	usual na	e II shall consist of the drugs and other substances, by whatever official name, me, chemical name, or brand name designated, listed in this section. Each drug assigned the Controlled Substances Code Number set forth opposite it.					
(b)	Narcotic Drugs: Substances, vegetable origin or chemical synthesis. Unless specifically excepted or unless listed in another schedule, 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 combination of extraction and chemical synthesis:						
	opia nalb naltr	am and opiate, and any salt, compound, derivative, or preparation of opium or te excluding apomorphine, thebaine-derived butorphanol, dextrorphan, uphine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, exone and samidorphan, and their respective salts, but including the owing:					
	(i)	Raw opium	9600*				
	(ii)	Opium extracts	9610*				
	(iii)	Opium fluid extracts	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*(9-2007)
		(xvi) Oxycodone	9143*
		(xvii)Oxymorphone	9652*
		(xviii) Thebaine	9333*
		(xix) Tapentadol	9780-(5-2009)
		(xx) Noroxymorphone	9668-(4-2021)
	(2)	Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in paragraph (b) (1) of this section, except that these substances shall not include the isoquinoline alkaloids of opium.*	
	(3)	Opium poppy and poppy straw.*	
	(4)	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;	
		(ii) [123I]ioflupane; or	
		(iii) [ <sup>18</sup> F]FP–CIT.	
	(5)	Concentrate or 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.*
(c)	foll wh	<b>Diates:</b> (Narcotic Drugs) Unless specifically excepted or unless in another schedulowing opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters, enever the existence of such isomers, esters, ethers, and salts is possible within the emical designations:	rs and ethers
	(1)	Alfentanil	9737-(2-1987)

(2) A	lphaprodine	9010*
(3) A	nileridine	9020*
(4) B	ezitramide	9800*
(5) B	sulk Dextropropoxyphene (non-dosage forms)	9273-(9-1981)
(6) C	arfentanil	9743-(9-1988)
(7) D	hydrocodeine	9120*
(8) D	iphenoxylate	9170*
(9) F	entanyl	9801*
(10)	Isomethadone	9226*
(11)	Levo-alphacetylmethadol (LAAM)	9648-(12-1993)
(12)	Levomethorphan	9210*
(13)	Levorphanol	9220*
(14)	Metazocine	9240*
(15)	Methadone	9250*
(16)	Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane	9254*
(17) c:	Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-arboxylic acid	9802*
(18)	Pethidine (Meperidine)	9230*
(19)	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine	9232*
(20)	Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate	9233*
(21)	Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid	9234*
(22)	Phenazocine	9715*
(23)	Piminodine	9730*
(24)	Racemethorphan	9732*
(25)	Racemorphan	9733*
(26)	Remifentanil	- 9739-(11-1996)
(27)	Sufentanil	9740-(9-1981)
(28)	Thiafentanil	9729-(4-2021)
(29)	Oliceridine	9245-(5-2022)
(30)	Tianeptine	(5-2022)

<sup>(</sup>d) **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:

(1) Amp	ohetamine, its salts, optical isomers, and salts of its optical isomers	1100*
(2) Met	namphetamine, its salts, isomers, and salts of its isomers	1105*
(3) Lisd	examfetamine	1205*(7-2007)
(4) Pher	nmetrazine and its salts	1631*
(5) Met	nylphenidate	1724*
compou depressa wheneve	cants: Unless specifically excepted or unless listed in another schedule, any and, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system, including its salts, isomers, and salts or the existence of such salts, isomers, and salts of isomers is possible within designation:	stances having a s of isomers
(1) Amo	barbital	2125*
(2) Glut	ethimide	2550-(2-1991)
(3) Pent	obarbital	2270*
(4) Pher	ncyclidine	7471*
(5) Seco	barbital	2315*
(f) Halluci	nogenic Substances:	
(1) Nab	ilone	- 7379-(11-1987)
	Other name(s) for nabilone: $(\pm)$ -trans-3- $(1,1$ -dimethylheptyl)-6,6a,7,8,10,10a exahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one].	-
U.S.	nabinol in an oral solution in a drug product approved for marketing by the Food and Drug Administration; [(-)-delta-9-trans-hydrocannabinol(delta-9-THC)]	7365-(7-2019)
	ate Precursor: Unless specifically excepted or unless listed in another sche, compound, mixture, or preparation which contains any quantity of the follows:	
(1) Imm	ediate precursor to Amphetamine and Methamphetamine:	
(i)	Phenylacetone	8501-(3-1980)
	Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl Ketone; methyl benzyl Ketone.	
(2) Imm	ediate precursor to Phencyclidine (PCP):	
(i)	1-phenylcyclohexylamine	7460*
(ii)	1-piperidinocyclohexanecarbonitrile (PCC)	8603*
(3) Imm	ediate precursor to Fentanyl:	
(i)	4-anilino-N-phenethylpiperidine (ANPP)	0222*(0.2010)

(ii) N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl) ----- 8366-(4-2021)

## **SCHEDULE III**

- (a) 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.
- (b) **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:
- (c) **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:
  - (1) Any compound, mixture, or preparation containing:

(i)	Amobarbital	2126*
(ii)	Secobarbital	2316*
(iii)	Pentobarbital	2271*
(iv)	Embutramide	2020*(9-2006)
	or any salt thereof and one or more other active medicinal ingredients which are not listed in any schedule.	

(2) Any suppository dosage form containing:

(i)	Amobarbital	2126*
(ii)	Secobarbital	2316*
(iii)	Pentobarbital	2271*

or any salt of any of these drugs and approved by the Food and Drug Administration for marketing only as a suppository.

	ny substance which contains any quantity of a derivative of barbituric acid or ny salt thereof	2100*
(4) C	hlorhexadol	2510*
is	ny drug product containing gamma hydroxybutyric acid, including its salts, omers, and salts of isomers, for which an application is approved under section of the Federal Food, Drug, and Cosmetic Act	- 2012-(2-2001)
(6) K	etamine. its salts, isomers, and salts of isomers	- 7285-(7-1999)
	Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
(7) L	ysergic acid	7300*
(8) L	ysergic acid amide	7310*
(9) M	lethyprylon	2575*
(10)	Sulfondiethylmethane	2600*
(11)	Sulfonethylmethane	2605*
(12)	Sulfonmethane	2610*
(13)	Tiletamine and zolazepam or any salt thereof	- 7295-(3-1988)
	Some trade or other name for a tiletamine- zolazepam 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-trimethyl-pyrazolo[3,4-e][1,4]diazepin-7(1H)-one; flupyrazapon.	
(14)	Perampanel	2261-(11-2013)
(15)	Xylazine and any material, compound, mixture, or preparation which contains by quantity of xylazine, including its salts, isomers, and salts of isomers	

- any quantity of xylazine, 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, except in the following uses:
  - Dispensing, prescribing, or administering, to an animal, a drug containing (i) xylazine that has been approved by the United States Secretary of Health and Human Services under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b);
  - (ii) Dispensing, prescribing, or administering xylazine to an animal that is permissible under section 512 (a)(4) of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(a)(4));
  - (iii) Possessing a drug containing xylazine, as described in this section (15), for animal use:

	(	A) by a need pharmacist of needsed vetermarian; or	
	(	B) Pursuant to a valid prescription from a licensed veterinarian;	
	(iv)	Possessing, manufacturing, distributing, or using xylazine as an active pharmaceutical ingredient for manufacturing an animal drug either:	
	(	(A) Approved under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b); or	
	(	B) Issued an investigation use exemption under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(j));	
	(v)	Manufacturing, distributing, or using a xylazine bulk chemical for pharmaceutical compounding by a licensed pharmacist or veterinarian; or	
	(vi)	Another use approved or permissible under the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 301, et seq.) or under 21 CFR Part 530, Subpart B.	
(d) N	Valorph	ine	9400*
(e) I	Narcot	ic drugs: Unless specifically excepted or unless listed in another schedule:	
(	narc	material, compound, mixture, or preparation containing any of the following otic drugs, or their salts calculated as the free anhydrous base or alkaloid, in ted quantities as set forth below:	
	(i)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*
	(ii)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9804*
	(iii)	Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9807*
	(iv)	Not more than 300 milligrams of ethylmorphine per 100 milliliters or not more than 15 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9808*
	(v)	Not more than 500 milligrams of opium per 100 milliliters or per 100 grams or not more than 25 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9809*
	(vi)	Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	
(2		material, compound, mixture, or preparation containing any of the following otic drugs or their salts, as set forth below:	

	(i)	Buprenorphine (10-02 Transfer)	9064-(6-1985)
	(ii	) Reserved	
(f)	comp salts,	olic Steroids: Unless specifically excepted or unless listed in another sche ound, mixture, or preparation containing any quantity of the following sub isomers, and salts of isomers whenever the existence of such salts of isome ecific chemical designation:	ostances, including its ers is possible within
	(1) B	oldenone;	(9-1991)
	(2) B	oldione;	(1-2010)
	(3) C	hlorotestosterone (4-chlortestosterone);	(9-1991)
	(4) C	lostebol;	(9-1991)
	(5) D	ehydrochlormethyltestosterone;	(9-1991)
	(6) D	esoxymethyltestosterone	(1-2010)
	(7) D	ihydrotestosterone (4-dihydrotestosterone);	(9-1991)
	(8) D	rostanolone;	(9-1991)
	(9) Et	hylestrenol;	(9-1991)
	(10)	Fluoxymesterone;	(9-1991)
	(11)	Formebulone (formebolone);	(9-1991)
	(12)	Mesterolone;	(9-1991)
	(13)	Methandienone, also known as Methandrostenolone;	(9-1991)
	(14)	Methandranone;	(9-1991)
	(15)	Methandriol;	(9-1991)
	(16)	Methenolone;	(9-1991)
	(17)	Methyltestosterone;	(9-1991)
	(18)	Mibolerone;	(9-1991)
	(19)	Nandrolone;	(9-1991)
	(20)	19-Nor-4,9(10)-Androstadienedione	(1-2010)
	(21)	Norethandrolone;	(9-1991)
	(22)	Oxandrolone;	(9-1991)
	(23)	Oxymesterone;	(9-1991)
	(24)	Oxymetholone;	(9-1991)
	(25)	Stanolone;	(9-1991)
	(26)	Stanozolol;	(9-1991)

(2	27)	Testolactone;	(9-1991)
(2	28)	Testosterone;	(9-1991)
(2	29)	Trenbolone	(9-1991)
(.	30)	Prostanozol	(8-2012)
(.	31)	Methasterone	(8-2012);
		and	
(.	32) pa	Any salt, ester, or isomer of a drug or substance described or list in this tragraph, if that salt, ester, or isomer promotes muscle growth	(9-1991)
		pt anabolic steroid products: Compounds, mixtures, or preparations that c d that have been exempted by the Secretary:	ontain an anabolic
		NDC Number	
(	1) A	ndro-Estro 90-4	0536-1605
(2	2) A	ndrogyn L.A	0456-1005
(.	3) C	omponent E-H in Process Pellets	Ivy Labs Inc.
(4	4) C	omponent E-H in Process Granulation	Ivy Labs Inc
(:	5) C	omponent TE-S in process Granulation	Ivy Labs Inc
(	6) C	omponent TE-S in process Pellets	Ivy Labs Inc
(	7) de	pANDROGYN	0456-1020
(8	8) D	epo-Testadiol	0009-0253
(9	9) D	EPO-T.E	52765-257
(	10)	depTESTROGEN	51698-257
(	11)	Duomone	52047-360
(	12)	DUO-SPAN II	0684-0102
(	13)	DURATESTRIN	43797-016
(	14)	Essian	Pharmaceutics
(	15)	Essian H.S	Pharmaceutics
(	16)	Esterified Estrogens & Methyltestosterone, USP (0.625 mg/1.25mg)	Interpharm
(	17)	Esterified Estrogens & Methyltestosterone, USP (1.25mg/2.5mg)	Interpharm
(	18)	Esterified Estrogens & Methyltestosterone (0.625mg/1.25mg) Tablet	ANDAPharm
(	19)	Esterified Estrogens & Methyltestosterone (1.25mg/2.5mg) Tablet	ANDAPharm
(2	20)	Estratest	0032-1026
(2	21)	Estratest HS	0032-1023
C	22)	Menogen	59243-570

(23)	Menogen HS	59243-560
(24)	Methyltestosterone & Esterified Estrogens (2.5mg/1.25Mg)	Lannett Co
(25)	Methyltestosterone & Esterified Estrogens (Half Strength) (1.25mg/0.625mg)	Lannett Co
(26)	PAN ESTRA TEST	0525-0175
(27)	Premarin with Methyltestosterone	0046-0879
(28)	Premarin with Methyltestosterone	0046-0878
(29)	Syntest D.S	66576-231
(30)	Stntest H.S	66576-230
(31)	Synovex H in process bulk pellets	Syntex Animal
(32)	Synovex H in process granulation	Syntex Animal
(33)	Synovex Plus in process granulation	Fort Dodge
(34)	Synovex Plus in process bulk pellets	Fort Dodge
(35)	TEST-ESTRO Cypionates	0536-9470
(36)	Testoderm with Adhesive 4mg/d	Alza Corp
(37)	Testoderm 4mg/d	17314-4608
(38)	Testoderm 6mg/d	17314-4609
(39)	Testoderm with Adhesive 6mg/d	17314-2836
(40)	Testoderm in process film	Alza Corp
(41)	Testoderm with Adhesive in process film	Alza Corp
(42)	Testosterone Cypionate/Estradiol Cypionate injection	54274-530
(43)	Testosterone Cypionate/Estradiol Cypionate injection	0182-3069
(44)	Testosterone Cyp 50 Estradiol Cyp 2	0814-7737
(45)	Testosterone Cypionate/Estradiol Cypionate injection	0364-6611
(46)	Testosterone Cypionate/Estradiol Cypionate injection	0402-0257
(47)	Testosterone Enanthate/Estradiol Valerate injection	0182-3073
(48)	Testosterone Enanthate/Estradiol Valerate injection	0364-6618
(49)	Testosterone Enanthate/Estradiol Valerate injection	0402-0360
(50)	Testosterone Ophthalmic Solution	Allergan
(51)	Tilapia Sex Reversal Feed (investigational)	Ranger, Inc

(h) Veterinary Anabolic Steroid Implant Products: Anabolic steroid products expressly intended for administration through implants in cattle or other nonhuman species exempted by the Secretary.

(1) Component E-H	021641-002
(2) Component E-H	01968327
(3) Component TE-S	021641-004
(4) Component T-H	0211641-006
(5) Component T-S	0211641-005
(6) F-TO	00093351
(7) Finaplix-H	12799-807-10
(8) Finaplix-S	12799-807-07
(9) Heifer-old	Boehringer
(10) Heifer-old	Ingelheim
(11) Heifer-old	Ivy Lab.
(12) Implus-H	0009-0434-01
(13) Implus-H	06-0434-01
01968327	
(14) Masculinizing Feed for Fish (Invesitigational)	Rangen,Inc.
(15) Revalor-G	12799-811
(16) Revalor-H	12799-810
(17) Revalor-S	12700 000
(11) 110.0000 2	12/99-809
(18) Synovex H	
	0856-3901
(18) Synovex H	0856-3901 Syntex

If veterinary products that are granted exempted status are subsequently distributed with the intent that they be used in humans, the distribution would be subject to the criminal sanctions of the CSA despite the drugs' exempted status.

## (i) Hallucinogenic substances:

(1) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a U.S. Food and Drug Administration approved drug product ----- 7369-(11-1987)

[Some other names for dronabinol: (6a R-trans)-6a,7,8, 10a-tetrahydro-6, 6, 9-trimethyl-3-pentyl-6H-dibenzo [b,d] phyran-1-ol, or (-)-delta 9-(trans)-tetrahydrocannabinol]

## SCHEDULE IV

- (a) 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.
- (b) <u>Narcotic drugs</u>: Unless specifically excepted or unless listed in another schedule, any material, 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 set forth below:
  - (1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit------ 9167\*
  - (2) Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-propionoxybutane) ------ 9278-(11-1987)
- (c) <u>Depressants</u>: 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) Alprazolam ----- 2882-(6-1982) (2) Barbital ------ 2145\* (3) Bromazepam ------ 2748-(1-1985) (4) Camazepam ------ 2749-(1-1985) (5) Chloral betaine ------2460\* (6) Chloral hydrate ----- 2465\* (7) Chlordiazepoxide ------ 2744\* (8) Clobazam ------ 2751-(1-1985) (9) Clonazepam ------ 2737\* (10) Clorazepate ------ 2768\* Clotiazepam ------ 2752-(1-1985) (11)Cloxazolam ------ 2753-(1-1985) (12)Delorazepam ------ 2754-(1-1985) (13)Diazepam ------ 2765\* Dichloralphenazone ------ 2467-(10-2002) (15)Estazolam ------ 2756-(1-1985) (16)Ethchlorvynol ------ 2540\* (17)Ethinamate ------ 2545\* (18)Ethyl loflazepate ------ 2758-(1-1985) (19)

(20)	Fludiazepam	2759-(1-1985)
(21)	Flunitrazepam	2763-(1-1985)
(22)	Flurazepam	2767*
(23)	Fospropofol	2138-(11-2009)
(24)	Halazepam	2762-(6-1982)
(25)	Haloxazolam	2771-(1-1985)
(26)	Ketazolam	2772-(1-1985)
(27)	Loprazolam	2773-(1-1985)
(28)	Lorazepam	2885*
(29)	Lormetazepam	2774-(1-1985)
(30)	Mebutamate	2800*
(31)	Medazepam	2836-(1-1985)
(32)	Meprobamate	2820*
(33)	Methohexital	2264*
(34)	Methylphenobarbital (mephorbarbital)	2250*
(35)	Midazolam	2884-(1-1985)
(36)	Nimetazepam	2837-(1-1985)
(37)	Nitrazepam	2834-(1-1985)
(38)	Nordiazepam	2838-(1-1985)
(39)	Oxazepam	2835*
(40)	Oxazolam	2839*
(41)	Paraldehyde	2585*
(42)	Petrichloral	2591*
(43)	Phenobarbital	2285*
(44)	Pinazepam	2883-(1-1985)
(45)	Prazepam	2764*
(46)	Quazepam	2881-(11-1986)
(47)	Temazepam	2925-(9-1981)
(48)	Tetrazepam	2886-(1-1985)
(49)	Triazolam	2887-(7-1983)
(50)	Zaleplon	2781-(9-1999)
(51)	Zolpidem	2783-(12-1993)
(52)	Zopiclone	2784-(1-2006)

(53)	Alfaxalone	2731-(2-2014)
(54)	Carisoprodol	8192-(4-1997)
(55)	Tramadol	9752-(8-2007)
(56)	Suvorexant	2223-(8-2014)
(57)	Brexanolone	2400-(4-2021)
(58)	Lemborexant	2245-(4-2021)
(59)	Remimazolam	2846-(5-2022)
(60)	Daridorexant	2410
(61)	Zuranolone	2420
stimul or geo isome	ound, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system, including its salts, isomers (whether operation), and salts of such isomers whenever the existence of such salts, isomers are is possible within the specific chemical designation:	ptical, position, , and salts of
	athine ((+)-Norpseudoephedrine)	
	iethylpropion	
` /	encamfamin	,
` ´	enproporex	· · · · · · · · · · · · · · · · · · ·
` /	orcaserin	` ′
` ′	azindol	` ` `
` /	efenorex	,
(8) M	odafinil	1680-(1-1999)
	emoline (including organometallic complexes and chelates thereof)	
(10)	Phentermine	1640*
(11)	Pipradrol	1750-(9-1981)
(12)	Serdexmethylphenidate	1729
(13)	Sibutramine	1675-(2-1998)
(14)	Solriamfetol	1650-(4-2021)
(15)	SPA ((-)-1-dimethylamino-1,2,diphenylethane)	1635-(9-1981)

(e) <u>Other substances</u>: 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 whether optical, position, or geometric), and salts of such isomers, whenever the existence of such salts, isomers, and salts of isomers is possible:

(1) Parters sing	0700 (4 1070)
(1) Pentazocine	,
(2) Butorphanol	,
(3) Nalbuphine	` ,
(4) Eluxadoline	9725-(4-2017)
SCHEDULE V	
(a) Schedule V shall consist of the drugs and other substances by whatever official nan usual name, chemical name, or brand name designated, listed in this section.	ne, common or
(b) <b>Narcotic Drugs:</b> Unless specifically excepted or unless listed in another schedule, compound, mixture or preparation containing any of the following narcotic drugs as set forth below.	
Reserved	
(c) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any comportant or preparation containing any of the following limited quantities of narcotic drugs of which shall include one or more nonnarcotic active medicinal ingredients in sufficient confer upon the compound, mixture, or preparation valuable medicinal qualities of possessed by the narcotic drug alone:	or salts thereof, ent proportion to
(1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams.	*
(2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams.	*
(3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams.	*
(4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit.	
(5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams	*
(6) Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.	*
(d) <u>Stimulants</u> : unless specifically excepted or unless listed in another schedule, any recompound, mixture, or preparation which contains any quantity of the following su stimulant effect on the central nervous system, including its salts, isomers and salts	bstances having
(1) Pyrovalerone	1485-(3-1988)
(2) Ephedrine:a -{-(Methylamino)ethyl}benzene-methanol;	(10-1995)

a-{-(methylamino) ethyl}benzyl alcohol; 2-methylamino-1-phenyl-1-propanol;

methylaminopropanol; a - hydroxy-b-methylaminopropylbenzene; a product

1-phenyl-1-hydroxy-2-methylaminopropane; 1-phenyl-2-

which occurs in the Chinese herb Ma Huang (Ephedra vulgaris, Ephedra sinica Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra spp.

(3) Phenylpropanolamine(	(7-	-2	0(	0:	5	
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(4) Pseudoephedrine ----- (7-2005)

Pursuant to Ark. Code Ann. § 5-64-212 as amended in 2005, this Schedule V classification shall NOT apply to any ephedrine, phenylpropanolamine, or pseudoephedrine in liquid, liquid capsule, or liquid gel capsule form. However, sales limits mandated by statute shall apply to all products with ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless of the dosage form.

(e) **Depressants:** 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, isomers, and salts of isomers:

(1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	- 2782-(1-2006)
(2) Lacosamide	- 2746-(5-2009)
(3) Brivaracetam	- 2710-(4-2017)
(4) Lasmiditan	- 2790-(4-2021)
(5) Cenobamate	- 2720-(4-2021)

(6) Ganaxolone ------ 2401

- (f) Other substances:
  - (1) None.

## **SCHEDULE VI \*\*\*\***

(a) In addition to any substance placed in Schedule VI by the Secretary of the Department of Health under § 5-64-214, any material, compound, mixture, or preparation, whether produced directly or indirectly from a substance of vegetable origin or independently by means of chemical synthesis, or by a combination of extraction and chemical synthesis, that contains any quantity of the following substances, or that contains any of their salts, isomers, and salts of isomers when the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation, is included in Schedule VI:

(1) Marijuana	*	k ×
(2) Tetrahydrocannabinols, unless the tetrahydrocannabinol is:	*	k ×

- (i) Contained in hemp-derived cannabidiol; -----(6-2020)
- (ii) Not more than three-tenths of one percent (0.3%) of delta-9 tetrahydrocannabinol in the hemp-derived cannabidiol on a dry weight

			d accuracy standards; and	(6-2020) ***
	(iii)		approved by the United States Food and Drug Administration for arketing as a medication;	(6-2020)
(3)	A sy	ntheti	c equivalent of:	
	(i)	The	substance contained in the Cannabis plant; or	**
	(ii)	The	substance contained in the resinous extractives of the genus Cannabis;	**
(4)	the p man extra salts	ified lant, a ufactured icts, in isom	botanically as Salvia divinorum, whether growing or not, the seeds of any extract from any part of the plant, and every compound, are, derivative, mixture, or preparation of the plant, its seeds, or its neluding salts, isomers, and salts of isomers when the existence of the ters, and salts of isomers is possible within the specific chemical in;	**
(5)	class spec Com num The	ses des ific ur pound erical	substances, derivatives, or their isomers in the chemical structural scribed below in subdivisions $(a)(5)(i)-(a)(5)(x)$ of this section and also neclassified substances in subdivision $(a)(5)(xi)$ of this section. ds of the structures described in this subdivision $(a)(5)$ , regardless of designation of atomic positions, are included in this subdivision $(a)(5)$ . etic substances, derivatives, or their isomers included in this subdivision	
	(i)	Tetra	ahydrocannabinols:	
	(	A) Te	trahydrocannabinols, including without limitation the following:	**
		a)	Delta-1 cis or trans tetrahydrocannabinol [other name(s): Delta-9 cis or trans tetrahydrocannabinol], and its optical isomers;	**
		b)	Delta-6 cis or trans tetrahydrocannabinol [other name(s): Delta-8 cis or trans tetrahydrocannabinol], and its optical isomers;	**
		c)	Delta- 3,4 cis or trans tetrahydrocannabinol [other name(s): Delta-6a,10a cis or trans tetrahydrocannabinol], and its optical isomers;	**
		d)	Delta-10 cis or trans tetrahydrocannabinol, and its optical isomers;	***
		e)	Delta-8 tetrahydrocannabinol acetate ester;	***
		f)	Delta-9 tetrahydrocannabinol acetate ester;	***
		g)	Delta-6a,10a, tetrahydrocannabinol acetate ester;	***
		h)	Delta-10 tetrahydrocannabinol acetate ester; and,	***
		i)	A product derived from industrial hemp that was produced as a result of a synthetic chemical process that converted the industrial hemp or a substance contained in industrial hemp into Delta-8, Delta-9, Delta 6a,10a, or Delta-10 tetrahydrocannabinol including their respective acetate esters.	***

	(B) Dronabinol in sesame oil and encapsulated in a soft gelatin capsule in a drug product approved by the United States Food and Drug Administration is not a tetrahydrocannabinol under this subdivision (a)(5)(i);	**
(ii)	Naphthoylindoles, or 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 and whether or not substituted in the naphthyl ring to any extent, including without limitation the following:	**
	(A) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**
	(B) JWH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	**
	(C) JWH-018, or 1-Propyl-3-(1-naphthoyl)indole;	
	(D) JWH-019, or 1-Hexyl-3-(1-naphthoyl)indole;	**
	(E) JWH-073, or 1-Butyl-3-(1-naphthoyl)indole;	**
	(F) JWH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(G) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(H) JWH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl) indole;	**
	(I) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole;	**
	(J) JWH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;	**
	(K) JWH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;	**
	(L) JWH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole;	**
	(M) AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole;	**
	(N) MAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1-naphthalenyl)-methanone;	**
	(O) EAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)-methanone; and	**
	(P) THJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl)methanone;	7024-(7-2019)
(iii)	Naphthylmethylindoles, or any compound structurally derived from an H-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 and whether or not substituted in the naphthyl ring to any extent, including without limitation the following:	**
	(A) JWH-175, or 1-Pentyl-1H-indol-3-vl-(1-naphthyl)methane; and	**

(	(B) JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1-naphthyl)methane;	**
(iv)	Naphthoylpyrroles, or 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 and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-307, or (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-ylmethanone;	**
(v)	Naphthylmethylindenes, or any compound structurally derived from 1-(1-napthylmethyl)indene with substitution at the 3-position of the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-176, or E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane;	**
(vi)	Phenylacetylindoles, or 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 and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
(	(A) JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;	**
	(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; and	
	(E) RCS-8, or 1-(2-cyclohexylethyl)-3-(2- methoxyphenylacetyl)indole;	
(vii)	Cyclohexylphenols, or 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 without limitation the following:	**
(	(A) CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]- phenol;	**
(	(B) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and	**
(	(C) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;	**

(vii	(ii) Benzoylindoles, or any compound structurally derived from 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 without limitation the following:	**
	(A) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;	
	(B) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;	**
	(C) WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;	**
	(D) AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole; and	**
	(E) RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1H-indol-3-yl)-methanone;	**
(ix)	Adamantoylindoles, or Adamantoylindazoles, including Adamantyl Carboxamide Indoles and Adamantyl Carboxamide Indazoles, or any compound structurally derived from 3-(1-adamantoyl) indole, 3-(1-adamantoyl) indazole, or 3-(2-adamantoyl)indole by substitution at a nitrogen atom of the indole or indazole ring with alkyl, haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl,cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole or indazole ring to any extent and whether or not substituted in the adamantly ring to any extent, including without limitation the following:	**
	(A) AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;	
	(B) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;	**
	(C) JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide [other name(s): 2NE1, moved in Schedule VI in 2020]	**
	(D) AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide;	**
	(E) 5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7049**
	(F) STS-135, or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide;	**
(x)	Tetramethylcyclopropylcarbonylindoles or any compound structurally derived from 3-(2,2,3,3-tetramethylcyclopropylcarbonyl) indole by substitution at the nitrogen atom of the indole ring with alkyl,haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl, cycloalkylethyl, (N-methylpiperidin-2-yl)methyl or 2-(4-morpholinyl)ethyl, whether or not	

		limitation the following:	**
	(A)	UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(B)	XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(C)	A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(D)	) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(E)	5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and	**
	(F)	A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(xi)	) L	Unclassified Synthetic Cannabinoids, including without limitation the following:	**
	(A)	CP 50556-1 hydrochloride, 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;	**
	(B)	) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;	**
	(C)	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;	**
	(D)	Dimethylheptylpyran or DMHP;	**
	(E)	WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone;	**
	(F)	URB597, or [3-(3-carbamoylphenyl)phenyl] N-Cyclohexylcarbamate;	**
	(G)	) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one;	**
	(H)	CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]-methanone;	**
		a) URB602, or cyclohexyl N-(3-phenylphenyl)carbamate;	**
	(I)	PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3-carboxylate;	**
	(J)	5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;	**
	(K)	BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate;	**
	(L)	NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide;	**
	(M	) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide;	**

	Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L- ine methyl ester	7033-(9-2018)
	MB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)-	
	Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-boxamido)-3,3-dimethylbutanoate;	7034-(11-2018)
	Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-boxamido)-3,3-dimethylbutanoate	7041-(11-2018)
	DMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H- indole-3-boxamido)-3,3-dimethylbutanoate;	7042-(11-2018)
	JB-AMB, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-boxamido)-3-methylbutanoate;	7021-(11-2018)
	DMB-FUBINACA, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-boxamido)-3,3-dimethylbutanoate;	7020-(11-2018)
	B-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-azole-3-caroboxamide;	7023-(7-2019)
	B-CHMINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-clohexylmethyl)-1H-indazole-3-carboxamide;	7031-(7-2019)
(W) 1-(c	MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-cyclohexylmethyl)-1H-indazole-3-carboxamide;	
(X) AE fluo	B-FUBINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-orobenzyl)-1H-indazole-3-carboxamide;	(9-2018)
	DB-PINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-indazole-3-carboxamide;	(9-2018)
	C-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-indazole-3-carboxamide	7083-(6-2020)
(AA) A	ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-luorobenzyl)-1H-indazole-3-carboxamide	- · 7010-(4-2021)
	4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-ndazole-3-carboxamido)-3,3-dimethylbutanoate;	7043-(4-2021)
	5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-luoropentyl)-1H-indazole-3-carboxamide	· 7025-(5-2022)
	4-CN-CUMYL-BUTINACA, or 1-(4-cyanobutyl)-N-(2-phenylpropan-yl)-1H-indazole-3-carboxamide	7089 –(5-2022)
	5F-CUMYL-P7AICA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-H-pyrrolo[2,3-b]pyridine-3-carboxamide	- 7085-(5-2022)
(FF) N	NM2201, or Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3- arboxylate	, ,

	(GG) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	7036
	(HH) FUB-144, or (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone;	7014
	(II) FUB-AKB48, or N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide;	7047
	(JJ) MDMB-4en-PINACA, or Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate;	
	(KK) CH-PIATA, or N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide;	
	(LL) ADB-BUTINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1H-indazole-3-carboxamide	7027
(6) A s	ynthetic substance, derivative, or its isomers with:	
(i)	Similar chemical structure to any substance described in subdivisions (a)(1)-(a)(5) of this section; or	**
(ii)	Similar pharmacological effects to any substance described in subdivisions (a)(1)-(a)(5) of this section.	**

- (b) However, except as provided under subsection (c) of this section, the Secretary shall not delete a controlled substance listed in this section from Schedule VI.
- (c) A prescription drug approved by the United States Food and Drug Administration under 21 U.S.C. § 355 is excluded from Schedule VI unless the secretary objects under § 5-64-201.
  - \*-Scheduled before April, 1979.
  - \*\*-Schedule VI is revised to conform to Act 329 of 2013.
  - \*\*\* Schedule VI is revised further to conform to Act 629 of 2023. Each substance added to the Controlled Substances List pursuant to Act 629 of 2023 shall have the following effective dates:
    - (a) For persons who are under twenty-one (21) years of age, the effective date shall be the effective date of Act 629 of 2023; and,
    - (b) For persons who are twenty-one (21) years of age or older, the effective date shall be August 1, 2023.
  - \*\*\*\* Pursuant to ongoing litigation, and a preliminary injunction against enforcing Act 629 of 2023, the changes made to the List of Controlled Substances pursuant to Act 629 of 2023 are not enforceable until a final order issued in the matter, Bio Gen, LLC, et al. v. Sarah Huckabee Sanders, et al., Case No. 4:23-CV-00718-BRW, Central Division, Eastern District of Arkansas, United States District Court.