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 10 a.m. on June 22, 2023, at the Arkansas Department of Health auditorium 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 www.healthy.arkansas.gov and at Pharmacy Services, Room #357, 3rd 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 June 22, 2023. Comments may be sent by email to shane.david@arkansas.gov or by mail to Shane David, Pharm.D., 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., Section Chief of Pharmacy Services 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., Section Chief Pharmacy Services Section
STATE OF ARKANSAS)	
COUNTY OF SALINE)	
I, Marci Middleton, do hereby certify that Shan me and signed the above referenced document.	e David, Pharm.D., well known to me, appeared befor
Sworn and subscribed to before me this	day of ,
	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

•	0	U	of controlled substanc the Board held in Littl	e Rock, Arkansas on th
day of	, ,	and after a Pu	ıblic Hearing on the	day of
,	, hel d in Little	Rock, Arkan	sas, at the State Depar	tment of Health Buildin
		· , - 	,	· · · · · · · · · · · · · · · · · · ·

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):

р	cetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-iperidinyl]-N-phenylacetamide)]	-9815-(2-1986)
	Acetylmethadol	
	Allylprodine	
(4) A	Alphacetylmethadol (except Levo-alphacetylmethadol (LAAM)	9603*
(5) A	Alphameprodine	9604*
(6) A	lphamethadol	9605*
p	Alpha-methylfentanyl [other name(s): (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-iperidyl] propronanilide; 1-(1-methyl-2-phenylethyl)-4(N-ropanilido)piperidine)]	9814-(6-1982)
n	Alpha-methylthiofentanyl(N-[1-methyl-2-(2thienyl)ethyl-4-piperidinyl]-N-henylpropanamide)	9832-(2-1986)
(9) B	Benzethidine	9606*
(10)	Betacetylmethadol	9607*
(11) p	Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-iperidinyl]-N-phenylpropanamide)]	9830-(2-1986)
(12)	Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-methyl-4-piperidinyl]-N-phenylpropamamide]	- 9831-(2-1986)
(13)	Betameprodine	9608*
(14)	Betamethadol	9609*
(15)	Betaprodine	9611*
(16)	Clonitazene	9612*
	Dextromoramide	06128
(17)	Dextromoramide	9013
(17)(18)	Diampromide	

(20)	Difenoxin 9	168*
(21)	Dimenoxadol 9	617*
(22)	Dimepheptanol 9	618*
(23)	Dimethylthiambutene 9	619*
(24)	Dioxaphetyl butyrate 9	621*
(25)	Dipipanone 9	622*
(26)	Ethylmethylthiambutene 9	623*
(27)	Etonitazene 9	624*
(28)	Etoxeridine 9	
(29)	Furethidine 9	
(30)	Hydroxypethidine 9	
(31)	Ketobemidone 9	
(32)	Levomoramide 9	
(33)	Levophenacylmorphan 9	631*
(34) pi	3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-iperidyl]-N-Phenylpropanamide)] 9813-(10-1	985)
(35) pl	3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-henylpropanamide) 9833-(2-1	986)
(36)	henylpropanamide) 9833-(2-1 Morpheridine 9	632*
(37)	MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)] 9661-(10-1	985)
(38)	Noracymethadol 9	
(39)	Norlevorphanol 9	
(40)	Normethadone 9	635*
(41)	Norpipanone 9	636*
(42) 4-	Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-piperindinyl]propananmide] 9812-(11-1	986)
(43)	PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] - 9663-(10-1	985)
(44)	Phenadoxone 9	637*
(45)	Phenampromide 9	638*
(46)	Phenomorphan 9	647*
(47)	Phenoperidine9	641*
(48)	Piritramide 9	642*
(49)	Proheptazine 9	643*

(50)	Properidine	9644*
(51)	Propiram	9649*
(52)	Racemoramide	9645*
(53)	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide	9835-(2-1986)
(54)	Tilidine	9750-(9-1981)
(55)	Trimeperidine	9646*
(56) pl	Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-henylacetamide]	-9821-(4-2017)
(57) pl	Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-henylbutyramide]	-9822-(4-2017)
	Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2-l)ethyl]piperidin-4-yl}-N-phenylpropionamide]	-9836-(4-2017)
(59) m	Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2-(4-nethylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	(4-2017)
	Valeryl fentanyl [other name(s): N-phenyl-N[1-(2-phenylethyl)-4-piperidinyl]-entanamide]	-9840-(4-2017)
	Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-henylfuran-2-carboxamide]	-9834-(4-2017)
(62) 4-	Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-piperidinyl]-propanamide]	-9827-(4-2017)
(63) pl	Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-henylethyl)piperidin-4-yl]acetamide]	-9838-(4-2017)
(64)	4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-henethylpiperidin-4-yl)butyramide]	
(65) pl	Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-henylethyl)-4-piperidinyl]-butanamide]	-9823-(4-2017)
(66)	Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide]	
(67) [(AH-7921 [other name(s): 3,4-dichloro- <i>N</i> -1dimethylamino)cyclohexylmethyl]benzamide]	-9551-(4-2017)
(68) cł	W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4-nlorophenyl)sulfonamide]	(4-2017)
	W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4-nlorophenyl)Sulfonamide]	(4-2017)
(70)	MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine]	9560-(4-2017)
(71) N	U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)-methylbenzamide]	9547-(4-2017)

	henylacrylamide]	9811-(6-2020)
(73) pl	4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-henethylpiperidin-4-yl)isobutyramide]	9824-(6-2020)
	Tetrahydrofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-henyltetrahydrofuran-2-carboxamide]	9843-(6-2020)
(75) pl	Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)- N-henylcyclopropanecarboxamide]	9845-(4-2021)
	Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-l)-N-phenylacetamide]	9825-(4-2021)
(77) pl	Ortho-fluorofentanyl [other name(s): N-(2-fluorophenyl)-N-(1-henethylpiperidin-4-yl)propionamide]	9816-(4-2021)
is 01 50	Fentanyl-related substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Fentanyl-related substance means any substance not therwise listed, and for which no exemption or approval is in effect under section 05 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that is tructurally 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;	
(i	Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(i	Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(i	Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; and/or	
(v	v) Replacement of the N-propionyl group by another acyl group.	
(79) pl	Crotonyl fentanyl [other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-henylbut-2-enamide]	9844-(5-2022)
	Cyclopentyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-henylcyclopentanecarboxamide]	9847-(5-2022)
	Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl)-N-(1-henethylpiperidin-4-yl)isobutyramide]	9826-(5-2022)
	Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-henethylpiperidin-4-yl)butyramide]	9837-(5-2022)
	Beta-methyl fentanyl [other name(s): N-phenyl-N-(1-(2-phenylpropyl) iperidin-4-yl)propionamide]	9856-(5-2022)
(84)	Beta'-phenyl fentanyl [other name: N-(1-phenethylpiperidin-4-yl)-N,3-iphenylpropanamide]	

` /	2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2-uorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	9855-(5-2022)
	4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) piperidin-	
4-	yl)-N-phenylacetamide]	9819-(5-2022)
	Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-nenethylpiperidin-4-yl)butyramide]	9846-(5-2022)
(88) ph	Ortho-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-menethylpiperidin-4-yl)acetamide]	-9848-(5-2022)
	Ortho-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2-ethylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9820-(5-2022)
(90) ph	Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1-menethylpiperidin-4-yl)propionamide]	9817-(5-2022)
	Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-nenylbenzamide]	9841-(5-2022)
(92) ph	Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-nenylthiophene-2-carboxamide]	9839-(5-2022)
(93) yl	Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4-)(phenyl)carbamate]	9851-(5-2022)
(94) ph	Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-nenethylpiperidin-4-yl)acrylamide]	9852-(5-2022)
(95) ph	Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-nenethylpiperidin-4-yl)isobutyramide]	9853-(5-2022)
	Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-nenethylpiperidin-4-yl)furan-2-carboxamide]	9854-(5-2022)
(97)	Isotonitazene	
(98)	Zipeprol	9873
sched the ex	m derivatives: (Narcotic Drugs) Unless specifically excepted or unless listed is ule, any of the following opium derivatives, its salts, isomers, and salts of isomeristence of such salts, isomers, and salts of isomers is possible within the specification:	rs whenever
(1) A	cetorphine	9319*
(2) A	cetyldihydrocodeine	9051*
(3) Be	enzylmorphine	9052*
(4) Co	odeine methylbromide	9070*
(5) Co	odeine-N-Oxide	9053*
(6) Cy	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)
mater hallue the ex desig	acinogenic substances: Unless specifically excepted or unless listed in another schedule, any rial, compound, mixture, or preparation, which contains any quantity of the following einogenic substance, or which contains any of its salts, isomers, and salts of isomers whenever eistence of such salts, isomers, and salts of isomers is possible within the specific chemical nation (for purposes of this paragraph only, the term "isomer" includes the optical, position and etric isomers):
(1) al	pha-ethyltryptamine 7249-(12-1993)
	Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET.
(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 aminoethane; alpha-desmethyl DOB; 2C-B, Nexus.

(4) 2,	5-dimethoxyamphetamine	7396*
	Some trade or other names: 2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA.	
(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- methylphenethylamine; paramethoxyamphetamine; PMA.	
(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 (methylenedioxy) phenethylamine, N-ethyl MDA; MDE; MDEA.	
(13)	N-hydroxy-3,4-methylenedioxyamphetamine	- 7402-(6-1990)
	Some trade or other names: N-hydroxy-alpha-methyl-3,4(methylenedioxy) phenethylamine; N-hydroxy MDA	
(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	
(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	
(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;	
2	5C-NBOMe	(5-2013)

	(30)	2-(4-10d0-2,3-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyljethanamine;	
	25	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	
	(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)
	(46)	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine; 25B-	
		BOMe	` /
	(47)	2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25B-BOH	(11_2018)
	(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH	
	(49)	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine;	· · · · · · · · · · · · · · · · · · ·
	` /	5E-NBOMe	
		2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 25H-	
		BOMe	(7-2019)
		2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]- phenol; 25C-BOH	(7_2019)
	(52)	2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25H-NBOH	,
	(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	
	(33)	Some trade or other names: Para-methoxymethamphetamine; PMMA	1243-(3-2022)
	(54)	2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one	7286
	(34)		/280
		Some trade or other names: Methoxetamine; MXE	
(e)	comp depre when	essants: Unless specifically excepted or unless listed in another schedule, any ound, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system, including its salts, isomers, and salt ever the existence of such salts, isomers, and salts of isomers is possible within ical designation:	ostances having a s of isomers
	(1) Pl	henazepam	(6-2012)
		amma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4-ydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate],

		known precursors and analogs. Precursors include but are not limited to: a-butyrolactone	-2010-(2-2001)
(3)	Meclo	qualone	2572*
(4)	Methad	qualone	2565*
(5)	Etizola	- m	(12-2014)
(6)	Clonaz	olam	(4-2021)
(7)	Flualpı	razolam	(4-2021)
(8)	Flubro	mazepam	(4-2021)
		mazolam	
Sti	mulant	s:	
(1)	compo substar	specifically excepted or unless listed in another schedule, any material, und, mixture, or preparation which contains any quantity of the following nees having a stimulant effect on the central nervous system, including its somers, and salts of isomers:	
	(i)	Cathinone	1235-(3-1988)
	(ii)	(±) CIS-4-Methylaminorex [(±)CIS-4,5-dihydro-4- methyl-5-phenyl-2-oxazolamine]	1590-(6-1990)
	(iii)	Fenethylline	1503-(9-1981)
	(iv)	N-Benzylpiperazine	7493-(1-2005)
		Some trade or other names: 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 (some other names: 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	1237-(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	

(f)

	(xi)	4,4'-Dimethylaminorex some other names: 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)
(2)	indirect chemic contain follow exister		
	(i)	4-Methylmethcathinone (Mephedrone)	
	(ii)	Methylenedioxypyrovalerone (MDPV)	
	(iii)	3,4-Methylenedioxy-N-methylcathinone (Methylone	7540-(3-2011)
	(iv)	4-Methoxymethcathinone	
	(v)	3-Fluoromethcathinone	
	(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)	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;	

	(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)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one (N-Ethylpentylone) 7543-((7-2019)
	(xix)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)7547-((4-2021)
	(xx)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone)	(4-2021)
	(xxi)	2-(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone; Alpha-Ethylaminohexanophenone)]	7246
	(xxii)	1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (Alpha-Pyrrolidinohexanophenone; Alpha-PHP)]	7544
	(xxiii)	2-(ethylamino)-1-(4-methylphenyl)pentan-1-one [other name(s): (4-Methyl-alpha-ethylaminopentiophenone; 4-MEAP)]	7245
	(xxiv)	1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (4'-Methyl-alpha-pyrrolidinohexiophenone; MPHP)]	
	(xxv)	1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha-Pyrrolidinoheptaphenone; PV8)]	7548
	(xxvi)	1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): (4'-Chloro-alpha-pyrrolidinovalerophenone; 4-chloro-alpha-PVP)]	7443
<u>SC</u>	HEDULE	II .	
(a)	usual name	If shall consist of the drugs and other substances, by whatever official name, comme, chemical name, or brand name designated, listed in this section. Each drug or substances the Controlled Substances Code Number set forth opposite it.	
(b)	or unless lindirectly	Drugs: Substances, vegetable origin or chemical synthesis. Unless specifically existed in another schedule, any of the following substances whether produced direct by extraction from substances of vegetable origin, or independently by means of chor by combination of extraction and chemical synthesis:	ly or
	opiate nalbup	and opiate, and any salt, compound, derivative, or preparation of opium or excluding apomorphine, thebaine-derived butorphanol, dextrorphan, thine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, tone and samidorphan, and their respective salts, but including the ing:	
	(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	
	(xiii)	Metopon	
	(xiv)	Morphine	
	(xv)	Oripavine	
	(xvi)	Oxycodone	9143*
	(xvii)	Oxymorphone	9652*
	(xviii)	Thebaine	
	(xix)	Tapentadol	9780-(5-2009)
	(xx)	Noroxymorphone	9668-(4-2021)
(2)	equival of this	It, compound, derivative, or preparation thereof which is chemically lent or identical with any of the substances referred to in paragraph (b) (1) section, except that these substances shall not include the isoquinoline ds of opium.*	
(3)	Opium	poppy and poppy straw.*	
(4)	leaves, derivat derivat	eaves (9040) and any salt, compound, derivative, or preparation of coca (including cocaine (9041) and ecgonine (9180) and their salts, isomers, ives and salts of isomers and derivatives), and any salt, compound, ive, or preparation thereof which is chemically equivalent or identical with 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)	[¹⁸ F]FP-CIT.	
` /	solid o	ntrate or poppy straw (the crude extract of poppy straw in either liquid, r powder form which contains the phenanthrene alkaloids of the opium	9670.*

(c) **Opiates:** (Narcotic Drugs) Unless specifically excepted or unless in another schedule, any of the following opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters and ethers whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designations:

(1) A	lfentanil	9737-(2-1987)
(2) A	lphaprodine	9010*
(3) A	nileridine	9020*
(4) B	ezitramide	9800*
(5) B	ulk Dextropropoxyphene (non-dosage forms)	9273-(9-1981)
(6) C	arfentanil	9743-(9-1988)
	hydrocodeine	
(8) D	iphenoxylate	9170*
(9) F	entanyl	9801*
(10)	Isomethadone	9226*
(11)	Levo-alphacetylmethadol (LAAM)	9648-(12-1993)
(12)	Levomethorphan	
(13)	Levorphanol	
(14)	Metazocine	9240*
(15)	Methadone	9250*
(16)	Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane	9254*
(17)	Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-	0002*
	Pethidine (Meperidine)	9802*
(18)		
(19)	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine	
(20)	Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate	
(21)	Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid	
(22)	Phenazocine	
(23)	Piminodine	
(24)	Racemethorphan	
(25)	Racemorphan	
(26)	Remifentanil	· · · · · · · · · · · · · · · · · · ·
(27)	Sufentanil	· · · · · · · · · · · · · · · · · · ·
(28)	Thiafentanil	· · · · · · · · · · · · · · · · · · ·
(29)	Oliceridine	
(30)	Tianeptine	(5-2022)

compour	nts: Unless specifically excepted or unless listed in another schedule, any mand, mixture, or preparation which contains any quantity of the following substate effect on the central nervous system:	
(1) Amp	hetamine, its salts, optical isomers, and salts of its optical isomers	1100*
(2) Meth	namphetamine, its salts, isomers, and salts of its isomers	1105*
(3) Lisdo	examefetamine	1205*(7-2007)
	metrazine and its salts	
(5) Meth	ylphenidate	1724*
compour depressa wheneve	ants: Unless specifically excepted or unless listed in another schedule, any mand, mixture, or preparation which contains any quantity of the following substitute effect on the central nervous system, including its salts, isomers, and salts of the existence of such salts, isomers, and salts of isomers is possible within the designation:	tances having a of isomers
	barbital	
	ethimide	
	obarbital	
	cyclidine	
(5) Seco	barbital	2315*
(f) Hallucir	nogenic Substances:	
(1) Nabi	lone	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.	abinol in an oral solution in a drug product approved for marketing by the Food and Drug Administration; [(-)-delta-9-trans-nydrocannabinol(delta-9-THC)]	-7365-(7-2019)
(g) Immedi	ate Precursor: Unless specifically excepted or unless listed in another sched compound, mixture, or preparation which contains any quantity of the follow	ule, any
(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) Immed	diate precursor to Fentanyl:	
	(i)	4-anilino-N-phenethylpiperidine (ANPP)	8333*(8-2010)
	(ii)	N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl)	8366-(4-2021)
<u>SC</u>	HEDULE	<u>: III</u>	
(a)	usual nam	III shall consist of the drugs and other substances, by whatever official nate, chemical name, or brand name designated, listed in this section. Each assigned the DEA Controlled Substances Code Number set forth opposite	drug or substance
(b)	compound stimulant or geomet	ts: Unless specifically excepted or unless listed in another schedule, any d, mixture, or preparation which contains any quantity of the following street on the central nervous system, including its salts, isomers (whether ic), and salts of such isomers whenever the existence of such salts, isomers possible within the specific chemical designation:	ubstances having a or optical, position,
	stimul prepar Sectio list for	e compounds, mixtures, or preparations in dosage unit form containing and lant substances listed in Schedule II which compounds, mixtures, or rations were listed on August 25, 1971, as excepted compounds under on 308.32, and any other drug of the quantitative composition shown in the those drugs or which is the same except that it contains a lesser quantity of substances	nat y of
		hetamine	
		phentermine	
	(4) Clorte	ermine	1647*
	(5) Phend	limetrazine	1615*
(c)	compound	nts : Unless specifically excepted or unless listed in another schedule, and, mixture, or preparation which contains any quantity of the following state effect on the central nervous system:	
	(1) Any c	ompound, 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 s	uppository dosage form containing:	
	(i)	Amaharhital	2126*

	(1	ii)	Secobarbital	2316*
	(1	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.	
		-	bstance which contains any quantity of a derivative of barbituric acid or t thereof	2100*
	(4) (Chlorh	exadol	2510*
	is 5	somer 605 of	rug product containing gamma hydroxybutyric acid, including its salts, s, and salts of isomers, for which an application is approved under section the Federal Food, Drug, and Cosmetic Act	2012-(2-2001)
	(6) K	Ketam:	ine. its salts, isomers, and salts of isomers	7285-(7-1999)
			ne other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-lohexanone.	
	(7) L	yserg	ic acid	7300*
		-	ic acid amide	
	(9) N		prylon	
	(10)		Condiethylmethane	
	(11)	Sulf	Conethylmethane	2605*
	(12)	Sulf	Conmethane	2610*
	(13)	Tile	tamine and zolazepam or any salt thereof	7295-(3-1988)
		Son Tela	ne trade or other name for a tiletamine- zolazepam combination product: azol.	
			ne trade or other names for tiletamine: -2(ethylamino)-2-(2-thienyl)-ohexanone.	
			ne trade or other names for zolazepam: -4(2-fluorophenyl)-6,8-dihydro-8,-trimethylpyrazolo-[3,4-e] [1,4,]-diazepin-7(1-H)-one. flupyrazapon.	
	(14)	Pera	ımpanel	2261-(11-2013)
(d)	Nalo	rphine	e	9400*
(e)	Nar	cotic	drugs: Unless specifically excepted or unless listed in another schedule:	
	n	arcoti	aterial, compound, mixture, or preparation containing any of the following c drugs, or their salts calculated as the free anhydrous base or alkaloid, in quantities as set forth below:	
	(1	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	

(i	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*
(i	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*
(i	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*
(1	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*
(1	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*
	any material, compound, mixture, or preparation containing any of the following arcotic drugs or their salts, as set forth below:	
(i		9064-(6-1985)
(i	i) Reserved	
comp salts, the sp	polic Steroids: Unless specifically excepted or unless listed in another schedule, and bound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is poperific chemical designation:	, including its ossible within 4000
comp salts, the sp	polic Steroids: Unless specifically excepted or unless listed in another schedule, and bound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is possible.	, including its ossible within 4000
comp salts, the sp (1) B	polic Steroids: Unless specifically excepted or unless listed in another schedule, and bound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is poperific chemical designation:	, including its ossible within 4000
comp salts, the sp (1) B (2) B	polic Steroids: Unless specifically excepted or unless listed in another schedule, and bound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is popecific chemical designation: Soldenone;	, including its ossible within 4000 (9-1991) (1-2010)
comp salts, the sp (1) B (2) B (3) C	polic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is propecific chemical designation: Soldenone;	, including its ossible within 4000 (9-1991) (1-2010) (9-1991)
comp salts, the sp (1) B (2) B (3) C (4) C	colic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is projectific chemical designation: Soldenone;	, including its ossible within 4000 (9-1991) (1-2010) (9-1991) (9-1991)
comp salts, the sp (1) B (2) B (3) C (4) C (5) D	polic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is properties chemical designation: Soldenone; Chlorotestosterone (4-chlortestosterone);	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
(1) E (2) E (3) C (4) C (5) D (6) D	colic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is properly chemical designation: Coldenone;	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (1-2010)
(1) E (2) E (3) C (4) C (5) E (6) E (7) E	colic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is projectific chemical designation: Coldenone;	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (1-2010) (9-1991) (1-2010) (9-1991)
(a) C (b) C (c) C	polic Steroids: Unless specifically excepted or unless listed in another schedule, and pound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is precific chemical designation: Boldenone; Chlorotestosterone (4-chlortestosterone); Chlorotestosterone (4-chlortestosterone; Chlorotestosterone (4-chlortestosterone); Chlorotestosterone (4-chlortestosterone); Chlorotestosterone (4-chlortestosterone); Chlorotestosterone (4-chlortestosterone);	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (1-2010) (9-1991) (9-1991) (9-1991) (9-1991)
(a) C (b) C (c) C	colic Steroids: Unless specifically excepted or unless listed in another schedule, and cound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is projectific chemical designation: Coldenone; Chlorotestosterone (4-chlortestosterone); Chehydrochlormethyltestosterone; Chehydrochlormethyltestosterone; Cheydrochlormethyltestosterone; Cheydrochlormethyltestosterone; Cheydrochlormethyltestosterone; Cheydrochlormethyltestosterone; Chlorotestosterone (4-dihydrotestosterone); Chlorotestosterone; Chlorotesto	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
comp salts, the sp (1) B (2) B (3) C (4) C (5) D (6) D (7) D (8) D (9) B	polic Steroids: Unless specifically excepted or unless listed in another schedule, and bound, mixture, or preparation containing any quantity of the following substances isomers, and salts of isomers whenever the existence of such salts of isomers is projectific chemical designation: Boldenone; Chlorotestosterone (4-chlortestosterone); Chlorotestosterone (4-chlortestosterone); Desoxymethyltestosterone; Dihydrotestosterone (4-dihydrotestosterone); Orostanolone; Chlylestrenol;	, including its ossible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (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;	
((20)	19-Nor-4,9(10)-Androstadienedione	
((21)	Norethandrolone;	(9-1991)
((22)	Oxandrolone;	(9-1991)
((23)	Oxymesterone;	` '
((24)	Oxymetholone;	(9-1991)
((25)	Stanolone;	
((26)	Stanozolol;	(9-1991)
((27)	Testolactone;	(9-1991)
((28)	Testosterone;	` '
((29)	Trenbolone	
((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 cod that have been exempted by the Secretary:	ntain an anabolic
			NDC Number
((1)	Andro-Estro 90-4	0536-1605
((2)	Androgyn L.A	0456-1005
((3) Co	omponent E-H in Process Pellets	Ivy Labs Inc.
((4) Co	omponent E-H in Process Granulation	Ivy Labs Inc
((5) Co	omponent TE-S in process Granulation	Ivy Labs Inc
((6) Co	omponent TE-S in process Pellets	Ivy Labs Inc
	` ′	pANDROGYN	•
((8) Do	epo-Testadiol	0009-0253

(9) D	DEPO-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
(20)	Estratest	0032-1026
(21)	Estratest HS	
(22)	Menogen	
(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	
(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	
(44)	Testosterone Cyp 50 Estradiol Cyp 2	
(45)	Testosterone Cypionate/Estradiol Cypionate injection	
(46)	Testosterone Cypionate/Estradiol Cypionate injection	
(47)	Testosterone Enanthate/Estradiol Valerate injection	
(48)	Testosterone Enanthate/Estradiol Valerate injection	
(49)	Testosterone Enanthate/Estradiol Valerate injection	
(50)	Testosterone Ophthalmic Solution	
(51)	Tilapia Sex Reversal Feed (investigational	_
` /	rinary Anabolic Steroid Implant Products: Anabolic steroid products exp nistration through implants in cattle or other nonhuman species exempted	pressly intended for
		NDC/DIN
` /	Component E-H	
(2)	Component E-H	01968327
	component TE-S	
(4) C	component T-H	0211641-006
	omponent T-S	
	-TO	
	inaplix-H	
	inaplix-S	
(9) H	leifer-old	Boehringer
(10)	Heifer-old	Ingelheim
(11)	Heifer-old	Ivy Lab.
(12)	Implus-H	
(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

	(20)	Synovex Plus	0856-3904
	(21)	Tilapia Sex Reversal Feed (investigational)	Zeigier Bros.
	that th	erinary products that are granted exempted status are subsequently distributed ney be used in humans, the distribution would be subject to the criminal sancte the drugs' exempted status.	
(i)	<u>Hallu</u>	cinogenic substances:	
		ronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule U.S. Food and Drug Administration approved drug product	
		[Some other names for dronabinol: (6a R-trans)-6a,7,8, 10a-tetrahydro-6, 6, trimethyl-3-pentyl-6H-dibenzo [b,d] phyran-1-ol, or (-)-delta 9-(trans)-tetrahydrocannabinol]	9-
<u>SC</u>	HEDU	JLE IV	
(a)	usual	lule IV shall consist of the drugs and other substances, by whatever official n name, chemical name or brand name designated, listed in this section. Each een assigned the DEA Controlled Substances Code Number set forth opposite	drug or substance
(b)	comp	otic drugs: Unless specifically excepted or unless listed in another schedule, ound, mixture, or preparation containing any of the following narcotic drugs, ated as the free anhydrous base or alkaloid, in limited quantities as set forth l	or their salts
	(1) N at	ot more than 1 milligram of difenoxin and not less than 25 micrograms of ropine sulfate per dosage unit	9167*
	(2) D	extro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-popionoxybutane)	9278-(11-1987)
(c)	comp	essants: Unless specifically excepted or unless listed in another schedule, an ound, mixture, or preparation which contains any quantity of the following staling its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation:	ubstances,
	(1) A	lprazolam	2882-(6-1982)
	(2) B	arbital	2145*
	(3) B	romazepam	2748-(1-1985)
	(4) C	amazepam	2749-(1-1985)
	(5) C	hloral betaine	2460*
	(6) C	hloral hydrate	2465*
	(7) C	hlordiazepoxide	2744*

(19) Synovex H ------ Syntex

(8) C	lobazam	2751-(1-1985)
(9) C	lonazepam	2737*
(10)	Clorazepate	2768*
(11)	Clotiazepam	2752-(1-1985)
(12)	Cloxazolam	2753-(1-1985)
(13)	Delorazepam	2754-(1-1985)
(14)	Diazepam	2765*
(15)	Dichloralphenazone	
(16)	Estazolam	2756-(1-1985)
(17)	Ethchlorvynol	2540*
(18)	Ethinamate	2545*
(19)	Ethyl loflazepate	2758-(1-1985)
(20)	Fludiazepam	
(21)	Flunitrazepam	2763-(1-1985)
(22)	Flurazepam	2767*
(23)	Fospropofol	
(24)	Halazepam	
(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	· /		
	(48)	Tetrazepam	2886-(1-1985)		
	(49)	Triazolam			
	(50)	Zaleplon			
	(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	,		
	(58)	Lemborexant	2245-(4-2021)		
	(59)	Remimazolam	2846-(5-2022)		
	(60)	Daridorexant	2410		
(d)	d) <u>Fenfluramine</u> : 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) Fe	enfluramine	1670*		
(e)	e) <u>Stimulants</u> : 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:				
	(1) C	athine ((+)-Norpseudeophedrine	1230-(3-1988)		
	(2) D	ethylpropion	1610*		
	(3) Fe	encamfamin	1760-(3-1988)		
	(4) Fe	nproporex	1575-(3-1988)		

	(5) Lord	caserin	- 1625-(6-2013)
	(6) Maz	rindol	- 1605-(6-1982)
	(7) Mef	enorex	- 1580-(3-1988)
	(8) Mod	dafinil	- 1680-(1-1999)
	(9) Pem	noline (including organometallic complexes and chelates thereof)	1530*
	(10)	Phentermine	1640*
	(11)	Pipradrol	- 1750-(9-1981)
	(12)	Serdexmethylphenidate	1729
	(13)	Sibutramine	
	(14)	Solriamfetol	
	(15)	SPA ((-)-1-dimethylamino-1,2,diphenylethane)	
(f)	f) Other substances: Unless specifically excepted or unless listed in another schedu compound, mixture, or preparation which contains any quantity of the following su including its salts; isomers whether optical, position, or geometric), and salts of su whenever the existence of such salts, isomers, and salts of isomers is possible:		stances,
	(1) Pen	tazocine	- 9709-(4-1979)
	(2) But	orphanol	- 9720-(4-1997)
		ouphine	
	(4) Elux	kadoline	- 9725-(4-2017)
86	HEDIH	EV	

SCHEDULE V

- (a) Schedule V shall consist of the drugs and other substances by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.
- (b) Narcotic Drugs: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation containing any of the following narcotic drugs and their salts, as set forth below.

Reserved

- (c) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following limited quantities of narcotic drugs or salts thereof, which shall include one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:
 - (1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams. -----*

	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 mate compound, mixture, or preparation which contains any quantity of the following substantial stimulant effect on the central nervous system, including its salts, isomers and salts of	ances having
	(1) Pyrovalerone	
	(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	, ,
	(3) Phenylpropanolamine(4) Pseudoephedrine	, ,
		ation shall aid capsule, or oroducts with
(e)	(4) Pseudoephedrine	ation shall aid capsule, or broducts with of the dosage
(e)	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 p ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless form. Depressants: Unless specifically exempted or excluded or unless listed in another schedule and the self-self-self-self-self-self-self-self-	ation shall aid capsule, or broducts with of the dosage hedule, any ving substances and salts of
(e)	(4) Pseudoephedrine	ation shall aid capsule, or broducts with of the dosage hedule, any ving substances and salts of
(e)	(4) Pseudoephedrine	ation shall aid capsule, or broducts with of the dosage hedule, any ving substances and salts of 2782-(1-2006) 2746-(5-2009)
(e)	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 p ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless form. Depressants: Unless specifically exempted or excluded or unless listed in another schedule and the selection of the follow having a depressant effect on the central nervous system, including its salts, isomers, a isomers: (1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	ation shall aid capsule, or broducts with of the dosage hedule, any ving substances and salts of 2782-(1-2006) 2746-(5-2009) 2710-(4-2017)
(e)	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. Depressants: Unless specifically exempted or excluded or unless listed in another sel 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: (1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	ation shall aid capsule, or broducts with of the dosage hedule, any ving substances and salts of 2782-(1-2006) 2746-(5-2009) 2710-(4-2017) 2790-(4-2021)

(f) Other substances:

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 -----** (2) Tetrahydrocannabinols, unless the tetrahydrocannabinol is: ----- ** Contained in hemp-derived cannabidiol; ----- (6-2020) (i) Not more than three-tenths of one percent (0.3%) of delta-9 (ii) tetrahydrocannabinol in the hemp-derived cannabidiol on a dry weight basis as verified by a nationally accredited laboratory for quality, purity and accuracy standards; and-----(6-2020) *** Not approved by the United States Food and Drug Administration for (iii) marketing as a medication; -----(6-2020) (3) A synthetic equivalent of: The substance contained in the Cannabis plant; or -----** (i) The substance contained in the resinous extractives of the genus Cannabis; (ii) _____** (4) Salvia divinorum or Salvinorin A, which includes all parts of the plant presently classified botanically as Salvia divinorum, whether growing or not, the seeds of the plant, any extract from any part of the plant, and every compound, manufacture, derivative, mixture, or preparation of the plant, its seeds, or its extracts, including salts, isomers, and salts of isomers when the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation; ------** (5) Synthetic substances, derivatives, or their isomers in the chemical structural classes described below in subdivisions (a)(5)(i)-(a)(5)(x) of this section and also specific unclassified substances in subdivision (a)(5)(xi) of this section. Compounds of the structures described in this subdivision (a)(5), regardless of numerical designation of atomic positions, are included in this subdivision (a)(5). The synthetic substances, derivatives, or their isomers included in this subdivision (a)(5) are: (i) Tetrahydrocannabinols:

(A) Tetrahydrocannabinols, including without limitation the following: -----**

8	a) Delta-1 cis or trans tetrahydrocannabinol [other name(s): Delta-9 cis or trans tetrahydrocannabinol], and its optical isomers;	**
ł	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	f)Delta-9 tetrahydrocannabinol acetate ester;	***
٤	g) Delta-6a,10a, tetrahydrocannabinol acetate ester;	
1	n) 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);	**
1 t ((i	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, eycloalkylmethyl, 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;	**

	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-(7-2
i i 1	Naphthylmethylindoles, or any compound structurally derived from an H-ndol-3-yl-(1-naphthyl) methane by substitution at the nitrogen atom of the ndole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, l-(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;
r t r s J	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-nethyl-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 WH-307, or (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-ylmethanone;
r t r s	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-nethyl-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 WH-176, or E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane;
I V I	Phenylacetylindoles, or any compound structurally derived from 3- ohenylacetylindole 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:

(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;	**
(В	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		
(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	

		limitation the following:	**
(A	A)	AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;	**
(E	3)	AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;	**
(0	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]	**
(I)	AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide;	**
(E	Ξ)	5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7049**
(F	7)	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:	**
(<i>A</i>	A)	UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(E	3)	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;	**
(I))	5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
(E	Ξ)	5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and	**
(F	7)	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	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;	**
(E	3)	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)	(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	e;
(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	
(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;	7034-(11-2018)
(Q)	5-Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate	7041-(11-2018)
(R)	MDMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H- indole-3-carboxamido)-3,3-dimethylbutanoate;	7042-(11-2018)
(S)	FUB-AMB, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate;	7021-(11-2018)
(T)	MDMB-FUBINACA, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	7020-(11-2018)
(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	(cyclohexylmethyl)-1H-indazole-3-carboxamide;	(11-2014)
	(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;	
	(Z)	5F-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	-7083-(6-2020)
	(Az	A) ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide	
	(BI	3) 4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	7043-(4-2021)
	(CC	C) 5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7025-(5-2022)
	(DI	2-yl)-1H-indazole-3-carboxamide	7089 –(5-2022)
	(EI	E) 5F-CUMYL-P7AICA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide	
	(FF	NM2201, or Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate	7221-(5-2022)
	(Go	G) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	7036
	(HI	H) 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;	
	(6) A syntl	netic 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)		except as provided under subsection (c) of this section, the Secretary shall necessition substance listed in this section from Schedule VI.	ot delete a
(c)	A prescrip	tion drug approved by the United States Food and Drug Administration und	ler 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.





Arkansas Department of Health

4815 West Markham Street ● Little Rock, Arkansas 72205-3867 ● Telephone (501) 661-2000
Governor Sarah Huckabee Sanders
Renee Mallory, RN, BSN, Interim Secretary of Health
Jennifer Dillaha, MD, Director

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.

- 1. Schedule I, (b), (11) is an item that has marked for clean-up.
- 2. AH-7921, MT-45 and U-47700 are Schedule 1 controlled substances. To follow DEA, controlled substance code numbers have been set forth opposite of each substance. Schedule I, (b), (67), Schedule I, (b), (70) and Schedule I, (b), (71).
- 3. Isotonitazene. 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. Schedule I, (b), (97).
- 4. Zipeprol. 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. Schedule I, (b), (98).
- 5. Methoxetamine, MXE. 2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one. The DEA has placed this hallucinogenic substance into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Schedule I, (d), (54).
- 6. Schedule I, (f), (1), (viii) is an item that has been marked for clean-up.
- 7. Amineptine. The DEA has placed this stimulant into Schedule I because it has not recognized medical use. This drug would be included as Schedule I to follow DEA. Schedule I, (f), (1), (xii).

- 8. Mesocarb. The DEA has placed this stimulant into Schedule I because it has not recognized medical use. This drug would be included as Schedule I to follow DEA with a subsequent numbering change. Schedule I, (f), (1), (xiii).
- 9. The items listed below are Schedule I controlled substances. To follow DEA controlled substance code numbers have been set forth opposite of each substance listed below. Schedule I, (f), (2), (vii) through (xvi).
 - 4-methyl-N-ethylcathinone (4-MEC)
 - 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)
 - Alpha-pyrrolidinopentiophenone (Alpha-PVP)
 - 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (Butylone) and this item has been marked for clean-up.
 - 2-(methylamino)-1-phenylpentan-1-one (Pentedrone)
 - 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone, bk-MBDP) and this item has been marked for clean-up.
 - 4-fluoro-N-methylcathinone (4-FMC; Flephedrone)
 - 3-fluoro-N-methylcathinone (3-FMC)
 - 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)
 - Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)
- 10. 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one, (N-Ethylpentylone) is a Schedule I controlled substance. This item has been marked for clean-up. Schedule I, (f), (2), (xviii).
- 11. 2-(ethylamino)-1-phenylhexan-1-one. N-Ethylhexedrone, Alpha-Ethylaminohexanophenone. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxi).
- 12. 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one. Alpha-Pyrrolidinohexanophenone, Alpha-PHP. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxii).
- 13. 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one. 4-Methyl-alphaethylaminopentiophenone, 4-MEAP. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This

- drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxiii).
- 14. 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one. 4'-Methyl-alpha-pyrrolidinohexiophenone, MPHP. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxiv).
- 15. 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one. Alpha-Pyrrolidinoheptaphenone, PV8. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxv).
- 16. 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one. 4'-Chloro-alpha-pyrrolidinovalerophenone, 4-chloro-alpha-PVP. The DEA has placed this synthetic cathinone into Schedule I because it has no recognized medical use. This drug would be included as Schedule I, to follow DEA. Schedule I, (f), (2), (xxvi).
- 17. [123I]ioflupane and [18F]FP–CIT have been identified as items specifically removed from schedules of the controlled substance Act. Language has been updated in Schedule II, (b), (4) to follow DEA. Schedule II, (b), (4) (i) through (iii). Updated language will reflect the following:
 - 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) $\int_{0}^{123} I \int_{0}^{123} I \int_{0}^{12$
 - (iii) $[^{18}F]FP$ –CIT.
- 18. Schedule II, (b), (5), is an item marked for clean-up.
- 19. Schedule II, (f), (1) is an item that has been marked for clean-up.

- 20. Schedule III, (c), (10), is an item marked for clean-up.
- 21. Schedule III, (f), specific prefatory language in anabolic steroids section is removed indicating *Items 1-28* and *(9-1991)* is removed and each specific substance that fits this designation will be identified by date separately, however; the controlled substance code number will remain in current position.
- 22. Schedule III, (f), (13) added language indicating Methandienone *also known as* Methandrostenolone.
- 23. Daridorexant. The FDA approved this drug for use in treatment of insomnia. This drug would be included as Schedule IV, to follow DEA. Schedule IV (c), (60).
- 24. Serdexmethylphenidate. The FDA approved this drug for use in treatment of attention deficit hyperactivity disorder. To follow DEA, this drug would be included as Schedule IV with subsequent numbering corrections to following this section. Schedule IV, (e), (12).
- 25. Lorcaserin and Solriamfetol are moved within Schedule IV, (e) to alphabetize the list. Schedule IV, (e), (5) and Schedule IV, (e), (14).
- 26. Ganaxolone. The FDA approved this drug for use in treatment of seizures associated with cyclin-dependent kinase-like 5 (CDKL5) deficiency disorder (CDD) in patients 2 years of age and older. This drug would be included as Schedule V to follow DEA. Schedule V, (e), (6).
- 27. Delta-1 cis or trans tetrahydrocannabinol, and its optical isomers is a Schedule VI controlled substance. The Arkansas State Crime Laboratory requested to include Delta-9 cis or trans tetrahydrocannabinol as an additional name of this substance to reflect an alternative chemical numbering system. Schedule VI, (a), (5), (i), (A), (a).
- 28. Delta-6 cis or trans tetrahydrocannabinol, and its optical isomers is a Schedule VI controlled substance. The Arkansas State Crime Laboratory requested to include Delta-8 cis or trans tetrahydrocannabinol as an additional name of this substance to reflect an alternative chemical numbering system. Schedule VI, (a), (5), (i), (A), (b).

- 29. Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers is a Schedule VI controlled substance that has been marked for clean-up. The Arkansas State Crime Laboratory requested to include Delta-6a,10a cis or trans tetrahydrocannabinol as an additional name of this substance to reflect an alternative chemical numbering system. Schedule VI, (a), (5), (i), (A), (c).
- 30. THJ-2201, AB-PINACA and AB-CHMINACA are Schedule VI controlled substances. To follow DEA, controlled substance code numbers have been set forth opposite of each substance. Schedule VI, (a), (5), (ii), (P), Schedule VI, (a), (5), (xi) (U) and Schedule VI, (a), (5), (xi), (V).
- 31. 5-Fluoro-MDMB-PICA and 5F-CUMYL-PINACA are Schedule VI controlled substances. Schedule VI, (a), (5), (xi), (Q) and Schedule VI, (a), (5), (xi), (Z). To follow DEA, a DEA Controlled Substance Code Number has been set forth opposite of each substance.
- 32. 5F-EDMB-PINACA. Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate. The DEA has scheduled this synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Schedule VI, (a), (5), (xi), (GG).
- 33. FUB-144. (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone. The DEA has scheduled this synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Schedule VI, (a), (5), (xi), (HH).
- 34. FUB-AKB48. N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide. The DEA has scheduled this synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Schedule VI, (a), (5), (xi), (II).
- 35. Documentation referencing Director is replaced to Secretary, as applicable, throughout the list in compliance with state law.
- 36. The controlled substance list has been updated to reflect dates of promulgation beside several listed substances.
- 37. As a general administrative update to the controlled substance list, clerical changes are reflected to the outline numbers within specific sections.

- 38. Pursuant to Act 629, concerning language in Schedule VI of the List of Controlled Substances, is amended., (a),(2), (ii). The language additions for Tetrahydrocannabinol are as followed:
 - (2) Tetrahydrocannabinols, unless the tetrahydrocannabinol is:
 - (ii) Not more than three-tenths of one percent (0.3%) of delta-9 tetrahydrocannabinol in the hemp-derived cannabidiol on a dry weight basis as verified by a nationally accredited laboratory for quality, purity and accuracy standards; and
- 39. Pursuant to Act 629, substances and language in Schedule VI of the List of Controlled Substances is amended, Page 28, (a), (5), (i), (A). Tetrahydrocannabinols shall include the following substance additions and language, Page 28, (a), (5), (A) (d) through (i):
 - (i) Tetrahydrocannabinols:
 - (A) Tetrahydrocannabinols, including without limitation the following: **
 - 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;
 - 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 the industrial hemp into Delta-8, Delta-9, Delta-6a, 10a or Delta-10 tetrahydrocannabinol included in their respective acetate esters.
- 40. 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.

QUESTIONNAIRE FOR FILING PROPOSED RULES WITH THE ARKANSAS LEGISLATIVE COUNCIL

	PARTMENT
	ARD/COMMISSION
BOA	ARD/COMMISSION DIRECTOR
COI	NTACT PERSON
ADI	DRESS
PHO	ONE NO. EMAIL_
NAI	ME OF PRESENTER(S) AT SUBCOMMITTEE MEETING
PRF	ESENTER EMAIL(S)
	INSTRUCTIONS
Que wha	rder to file a proposed rule for legislative review and approval, please submit this Legislative estionnaire and Financial Impact Statement, and attach (1) a summary of the rule, describing at 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 R	he rule is being filed for permanent promulgation, please email these items to the attention tebecca Miller-Rice, miller-ricer@blr.arkansas.gov, for submission to the Administrative es Subcommittee.
Dire	ne rule is being filed for emergency promulgation, please email these items to the attention of ector Marty Garrity, garritym@blr.arkansas.gov , for submission to the Executive committee.
Plea	se answer each question completely using layman terms.
***	************************
1.	What is the official title of this rule?
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.	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.

8.	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).
9.	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).
10.	What is the reason for this proposed rule? Why is it necessary?

11.	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).
12.	Will a public hearing be held on this proposed rule? Yes No
	If yes, please complete the following:
	Date:
	Time:
	Place:
	Please be sure to advise Bureau Staff if this information changes for any reason.
13.	On what date does the public comment period expire for the permanent promulgation of the rule? Please provide the specific date.
14.	What is the proposed effective date for this rule?
15.	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.
16.	Please attach proof of filing the rule with the Secretary of State, as required by Ark. Code Ann. § 25-15-204(e)(1)(A).
17.	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.
18.	Is the rule expected to be controversial? Yes No If yes, please explain.

FINANCIAL IMPACT STATEMENT

PLEASE ANSWER ALL QUESTIONS COMPLETELY.

	PARTMENT
	ARD/COMMISSION
PER	SON COMPLETING THIS STATEMENT
TEL.	EPHONE NOEMAIL
email	omply with Ark. Code Ann. § 25-15-204(e), please complete the Financial Impact Statement and lit with the questionnaire, summary, markup and clean copy of the rule, and other documents. se 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?

	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 st	tate 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, amen rule, and explain how they are affected Current Fiscal Year	al year to any private individual, private entity, or private aded, or repealed rule? Please identify those subject to the l. Next Fiscal Year \$
í.	business subject to the proposed, amen	ided, or repealed rule? Please identify those subject to the l.
	business subject to the proposed, amenule, and explain how they are affected Current Fiscal Year \$	Next Fiscal Year Next Fiscal Year Subject to the labeled subject t
5.	business subject to the proposed, amenule, and explain how they are affected Current Fiscal Year \$	nded, or repealed rule? Please identify those subject to the l. Next Fiscal Year \$

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.

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 <u>Director Secretary</u> 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 <u>Director Secretary</u> 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 <u>Director Secretary</u> of the Arkansas Department of Health. A "*" denotes the substance was scheduled prior to April, 1979.

I, Shane David, Pharm.D., Section Chief of Pharmacy Services 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., Section Chief Pharmacy Services Section
STATE OF ARKANSAS)
COUNTY OF SALINE))
I, Marci Middleton, do hereby me and signed the above refer	y certify that Shane David, Pharm.D., well known to me, appeared before renced document.
Sworn and subscribed	to before me this day of ,
	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 Stat. 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

Rock, Arkansas, at th	a State Departmen	_ day of nt of Health Ruildir
 NUCK, AI Kallsas, at th	le State Departmen	it vi iicaitii Duiiuli

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):

	Acetyl-alpha-methylfentanyl <u>[other name(s):</u> (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	-9815-(2- <u>19</u> 86)
	Acetylmethadol	
(3)	Allylprodine	9602*
(4)	Alphacetylmethadol (except Levo-alphacetylmethadol (LAAM)	9603*
	Alphameprodine	
(6)	Alphamethadol	9605*
(7)	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- <u>19</u> 82)
(8)	Alpha-methylthiofentanyl(N-[1-methyl-2-(2thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	9832-(2- <u>19</u> 86)
(9)		
(10) Betacetylmethadol	9607*
(11	Beta-hydroxyfentanyl <u>[other name(s):</u> (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl []-N-phenylpropanamide)]	9830-(2- <u>19</u> 86)
(12	Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide]	
(13) Betameprodine	9608*
(14) Betamethadol	9609*
(15) Betaprodine	9611*
(16) Clonitazene	9612*
(17) Dextromoramide	9613*
(18) Diampromide	9615*
(19) Diethylthiambutene	9616*

(20)	Difenoxin	9168*
(21)	Dimenoxadol	9617*
(22)	Dimepheptanol	9618*
(23)	Dimethylthiambutene	9619*
(24)	Dioxaphetyl butyrate	9621*
(25)	Dipipanone	9622*
(26)	Ethylmethylthiambutene	
(27)	Etonitazene	
(28)	Etoxeridine	9625*
(29)	Furethidine	
(30)	Hydroxypethidine	
(31)	Ketobemidone	9628*
(32)	Levomoramide	
(33)	Levophenacylmorphan	9631*
(34) pi	3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-iperidyl]-N-Phenylpropanamide)]	9813-(10- <u>19</u> 85)
(35)	3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-henylpropanamide)	9833-(2-1986)
(36)	Morpheridine	9632*
(37)	MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)]	
(38)	Noracymethadol	
(39)	Norlevorphanol	
(40)	Normethadone	
(41)	Norpipanone	9636*
(42)	Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-piperindinyl]propananmide]	9812-(11- <u>19</u> 86)
(43)	PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine]	9663-(10- <u>19</u> 85)
(44)	Phenadoxone	9637*
(45)	Phenampromide	9638*
(46)	Phenomorphan	9647*
(47)	Phenoperidine	9641*
(48)	Piritramide	9642*
(49)	Proheptazine	9643*

(50)	Properidine	9644*
(51)	Propiram	9649*
(52)	Racemoramide	9645*
(53)	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide	9835-(2- <u>19</u> 86)
(54)	Tilidine	9750-(9- <u>19</u> 81)
(55)	Trimeperidine	9646*
(56) pł	Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-nenylacetamide]	-9821 <u>-(4-2017)</u>
(57) pl	Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-nenylbutyramide]	-9822 <u>-(4-2017)</u>
	Beta-hydroxythiofentanyl <u>[other name(s):</u> N-{1-[2-hydroxy-2-(thiophen-2-)ethyl]piperidin-4-yl}-N-phenylpropionamide]	-9836 <u>-(4-2017)</u>
	Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2-(4-ethylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	<u>(4-2017)</u>
	Valeryl fentanyl <u>[other name(s):</u> N-phenyl-N[1-(2-phenylethyl)-4-piperidinyl]-entanamide]	-9840 <u>-(4-2017)</u>
	Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-nenylfuran-2-carboxamide]	-9834 <u>-(4-2017)</u>
(62) 4-	Isobutyryl fentanyl <u>[other name(s):</u> 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-piperidinyl]-propanamide]	-9827 <u>-(4-2017)</u>
(63) pł	Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-menylethyl)piperidin-4-yl]acetamide]	-9838 <u>-(4-2017)</u>
(64) pł	4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-menethylpiperidin-4-yl)butyramide]	<u>(4-2017)</u>
	Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-nenylethyl)-4-piperidinyl]-butanamide]	-9823 <u>-(4-2017)</u>
(66)	Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide]	<u>(4-2017)</u>
(67) [(AH-7921 [other name(s): 3,4-dichloro- <i>N</i> -1dimethylamino)cyclohexylmethyl]benzamide]	- <u>9551-(4-2017)</u>
	W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4-nlorophenyl)sulfonamide]	<u>(4-2017)</u>
(69) cł	W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4-nlorophenyl)Sulfonamide]	<u>(4-2017)</u>
(70)	MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine]	9560-(4-2017)
(71) N	U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)-methylbenzamide]	9547-(4-2017)

phenylacrylamide]	9811-(6-2020)
(73) 4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	
(74) Tetrahydrofuranyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide]	
(75) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)- N-phenylcyclopropanecarboxamide]	9845 <u>-(4-2021)</u>
(76) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	
(77) Ortho-fluorofentanyl <u>[other name(s):</u> N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9816 <u>-(4-2021)</u>
(78) Fentanyl-related substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Fentanyl-related substance means any substance not otherwise listed, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that is structurally related to fentanyl by one or more of the following modifications:	
(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; and/or	
(v) Replacement of the N-propionyl group by another acyl group.	
(79) Crotonyl fentanyl <u>[other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide]</u>	9844 <u>-(5-2022)</u>
(80) Cyclopentyl fentanyl <u>[other name(s):</u> N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide]	9847 <u>-(5-2022)</u>
(81) Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9826 <u>-(5-2022)</u>
(82) Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	9837 <u>-(5-2022)</u>
(83) Beta-methyl fentanyl <u>[other name(s):</u> N-phenyl-N-(1-(2-phenylpropyl) piperidin-4-yl)propionamide]	9856 <u>-(5-2022)</u>
(84) Beta'-phenyl fentanyl <u>[other name:</u> N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide]	

	(85) 2'-Fluoro ortho-fluorofentanyl <u>lother name(s):</u> N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	9855 <u>-(5-2022)</u>
	(86) 4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) piperidin-4-yl)-N-phenylacetamide]	9819-(5-2022)
	(87) Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	
	(88) Ortho-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	-9848 <u>-(5-2022)</u>
	(89) Ortho-methyl methoxyacetyl fentanyl <u>[other name(s):</u> 2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9820 <u>-(5-2022)</u>
	(90) Para-methylfentanyl <u>[other name(s):</u> N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9817 <u>-(5-2022)</u>
	(91) Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide]	
	(92) Thiofuranyl fentanyl <u>[other name(s):</u> N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide]	9839 <u>-(5-2022)</u>
	(93) Fentanyl carbamate <u>[other name(s):</u> ethyl(1-phenethylpiperidin-4-yl)(phenyl)carbamate]	9851 <u>-(5-2022)</u>
	(94) Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide]	9852 <u>-(5-2022)</u>
	(95) Ortho-fluoroisobutyryl fentanyl <u>[other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]</u>	9853 <u>-(5-2022)</u>
	(96) Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	9854 <u>-(5-2022)</u>
	(97) Isotonitazene	
	(97)(98)Zipeprol	
(c)	Opium derivatives: -(Narcotic Drugs) Unless specifically excepted or unless listed schedule, any of the following opium derivatives, its salts, isomers, and salts of isomer the existence of such salts, isomers, and salts of isomers is possible within the specific designation:	rs whenever
	(1) Acetorphine	9319*
	(2) Acetyldihydrocodeine	9051*
	(3) Benzylmorphine	9052*
	(4) Codeine methylbromide	9070*
	(5) Codeine-N-Oxide	
	(6) Cyprenorphine	9054*

(7) I	Desomorphine	9055*
(8) I	Dihydromorphine	9145*
(9) I	Orotebanol	9335*
(10)	Etorphine (except hydrochloride salt)	9056*
(11)	Heroin	9200*
(12)	Hydromorphinol	9301*
(13)	Methyldesorphine	
(14)	Methyldihydromorphine	
(15)	Morphine methylbromide	9305*
(16)	Morphine methylsulfonate	9306*
(17)	Morphine-N-Oxide	9307*
(18)	Myrophine	
(19)	Nicocodeine	
(20)	Nicomorphine	9312*
(21)	Normorphine	9313*
(22)	Pholcodine	9314*
(23)	Thebacon	9315*
(24)	Mitragynine	· <u>(11-2015)</u>
(25)	7-Hydroxymitragynine	<u>(11-2015)</u>
mate hallu the e desig	ducinogenic substances: Unless specifically excepted or unless listed in another serial, compound, mixture, or preparation, which contains any quantity of the following constance, or which contains any of its salts, isomers, and salts of isomers existence of such salts, isomers, and salts of isomers is possible within the specific gnation (for purposes of this paragraph only, the term "isomer" includes the optical metric isomers):	wing ers whenever chemical
(1) a	alpha-ethyltryptamine	7249-(12- <u>19</u> 93)
	Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET.	
(2) 4	1-bromo-2,5-dimethoxy-amphetamine	7391*
	Some trade or other names: 4-bromo-2,5-dimethoxy-alphamethylphenethylamine; 4-bromo-2,5-DMA.	
(3) 4	1-bromo-2,5-dimethoxyphenethylamine	7392-(8- <u>19</u> 95)
	Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1 aminoethane; alpha-desmethyl DOB; 2C-B, Nexus.	

(4) 2,	,5-dimethoxyamphetamine	7396*
	Some trade or other names: 2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA.	
(5) 2,	,5-dimethoxy-4-ethylamphetamine	· 7399-(3- <u>19</u> 88)
	Some trade or other names: DOET.	
(6) 2,	,5-dimethoxy-4-(n)-propylthiophenethylamine	· 7348-(1- <u>20</u> 05)
	Some trade or other names: 2C-T-7.	
(7) 4	-methoxyamphetamine	7411 *
	Some trade or other names: 4-methoxy-alpha- methylphenethylamine; paramethoxyamphetamine; PMA.	
(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- <u>19</u> 85)
	Some trade or other names: MDMA)	
(12)	3,4-methylenedioxy-N-ethylamphetamine	7404-(6- <u>19</u> 90)
	Some trade or other names: N- ethy-alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA; MDE; MDEA.	
(13)	N-hydroxy-3,4-methylenedioxyamphetamine	7402-(6- <u>19</u> 90)
	Some trade or other names: N-hydroxy-alpha-methyl-3,4(methylenedioxy) phenethylamine; N-hydroxy MDA	
(14)	3,4,5-trimethoxy amphetamine	7390*
(15)	5-methoxy-n,n-dimethyltryptamine 5-MeO-DMT	7431*(1- <u>20</u> 11)
(16)	alpha-methyltryptamine	7432-(7- <u>20</u> 05)
	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- <u>20</u> 05)
	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	
(24)	Parahexyl	7374-(7- <u>19</u> 83)
	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- <u>19</u> 89)
	Some other trade or other names: TCPy.	
(34)	N,N-Diallyl-5-Methoxytryptamine;	<u>(6-2012</u>)
	Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT	
(35)	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine;	
2	5C-NBOMe	(5-2013 ⁾

		2-(4-10do-2,3-dimethoxyphenyi)-N-[(2-methoxyphenyi) methyljethanamne;	
	25	5I-NBOMe	<u>(8-2013)</u>
	(37)	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E	
	(38)	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D	7508 <u>-(11-2013)</u>
	(39)	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C	- 7519 <u>-(11-2013)</u>
	(40)	2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I	- 7518 <u>-(11-2013)</u>
	(41)	2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-2	- 7385 <u>-(11-2013)</u>
	(42)	2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-4	- 7532 <u>-(11-2013)</u>
	(43)	2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H	- 7517 <u>-(11-2013)</u>
	(44)	2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N	- 7521 <u>-(11-2013)</u>
	(45)	2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine; 2C-P	- 7524 <u>-(11-2013)</u>
	(46)	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine; 25B-BOMe	(9-2018)
		2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25B-	<u>(7-2016)</u>
	N	BOH	<u>(11-2018)</u>
	(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH	[<u>(11-2018)</u>
	(49)	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine;	(7.2010)
		5E-NBOMe	<u>(7-2019)</u>
		2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 25H-BOMe	(7-2019)
		2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]- phenol; 25C-	()
		BOH	<u>(7-2019)</u>
	(52)	2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25H-NBOH	<u>(7-2019)</u>
	(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	1245 <u>-(5-2022)</u>
		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	
. ,	comp depre when	essants: Unless specifically excepted or unless listed in another schedule, any ound, mixture, or preparation which contains any quantity of the following substant 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 ical designation:	stances having a s of isomers
	(1) Pl	nenazepam	<u>(6-2012)</u>
		amma-hydroxybutyric acid [other name(s): (some other names include GHB; amma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxydutanoic acid; sodium	

			e; sodium oxybutyrate), , and its known precursors and analogs. Precursors but are not limited to: gamma-butyrolactone	
	(3)		oqualone	
	` ′		qualone	
	` ′		um	
	` /		zolam	
	` ′		razolam	
			mazepam	
			mazolam	
(f)	Sti	mulant	s:	
	(1)	compo	ess specifically excepted or unless listed in another schedule, any material, bund, mixture, or preparation which contains any quantity of the following nees having a stimulant effect on the central nervous system, including its somers, and salts of isomers:	
		(i)	(1) Cathinone	1235-(3- <u>19</u> 88)
		(ii)	(2) (±) CIS-4-Methylaminorex [(±)CIS-4,5-dihydro-4- methyl-5-phenyl-2-oxazolamine]	1590-(6- <u>19</u> 90)
		(iii)	(3) Fenethylline	1503-(9- <u>19</u> 81)
		(iv)	(4) N-Benzylpiperazine	7493-(1- <u>20</u> 05)
			Some trade or other names: BZP, 1-Benzylpiperazine	
		(v)	(5) N-ethylamphetamine	1475-(6- <u>19</u> 82)
		(vi)	(6) N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers	9818-(2- <u>19</u> 86)
		(vii)	(7)-N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropan-mide (thenylfentanyl), its optical isomers, salts_and salts of isomers	9834-(2- <u>19</u> 86)
		(viii)	(8) N, N, N, N. Dimethylamphetamine <u>[other name(s): (some other names: N, N, Alpha-trimethylbenzeneethanamine; N, N, Alpha-trimethylphenethylamine),]</u> , its salts, optical isomers, and salts of optical isomers	1480-(2- <u>19</u> 89)
		(ix)	(9)—Methcathinone (some other names: 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 -	1237-(12- <u>19</u> 93)
		(x)	(10)—Aminorex [other name(s): (some other names: aminoraphen, 2-amino-5 phenyl-2-oxazoline, or 4,5 dihyrdo-5-phenyl-2-oxazolamine], its salts, optical isomers, and salts of optical isomers	1585-(12- <u>19</u> 93)

	(X1)	dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine, or 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine	1595 <u>-(5-2022)</u>
	(xii)	Amineptine	1219
	(xi)(xi	ii) Mesocarb	
	(xii)(x	iv)(12) Methyl-N-ethylcathinone (MEC)	<u>(6-2014)</u>
(2)	means synthe any of the exi the spe	Any material, compound, mixture, or preparation, whether produced y or indirectly from a substance of vegetable origin or independently by of chemical synthesis or by a combination of extraction and chemical sis, that contains any quantity of the following substances, or that contains the following substances' analogs, salts, isomers, and salts of isomers when stence of the analogs, salts, isomers, and salts of isomers is possible within the crific chemical designation, with the following chemical structure is ed in Schedule I:	
	(i)	(1) 4-Methylmethcathinone (Mephedrone)	
	(ii)	(2) Methylenedioxypyrovalerone (MDPV)	
	(iii)	(3)—3,4-Methylenedioxy-N-methylcathinone (Methylone	7540-(3- <u>20</u> 11)
	(iv)	(4) 4-Methoxymethcathinone	
	(v)	(5) 3-Fluoromethcathinone	(3- <u>20</u> 11)
	(vi)	(6) 4-Fluoromethcathinone-or	(3- <u>20</u> 11)
	(vii)	(7)—1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one_(Butylone) (7541-(11-2014)
	(viii)	(8)—Alpha-Pyrrolidinopentiophenone (Alpha-PVP)	7545-(11-2015)
	(ix)	(9) 4-methyl-N-ethylcathinone (4-MEC)	1249(9-2018)
	(x)	(10) 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)	-7498-(9-2018)
	(xi)	(11)-2-(methylamino)-1-phenylpentan-1-one (Pentedrone)	-1246-(9-2018)
	(xii)	(12) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)	- <u>7542-(9-2018)</u>
	(xiii)	(13) 4-fluoro-N-methylcathinone (4-FMC, Flephedrone)	-1238-(9-2018)
	(xiv)	(14)-3-fluoro-N-methylcathinone (3-FMC)	-1233-(9-2018)
	(xv)	(15) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)	-1258-(9-2018)
	(xvi)	(16) Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)	- <u>7546-(9-2018)</u>
	(xvii)	(17) 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:	

	(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)	(18) N-Ethylpentylone, or 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan- 1-one (N-Ethylpentylone)7543-(**)	<u>7-2019)</u>
	(xix)	(19)1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)7547-(4)	<u>4-2021)</u>
	(xx)	-(20)1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone -(Eutylone)(20)	<u>4-2021)</u>
	(xxi)	2-(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone; Alpha-Ethylaminohexanophenone)]	<u> 7246</u>
	(xxii)	1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (Alpha-Pyrrolidinohexanophenone; Alpha-PHP)]	7544
	(xxiii)	2-(ethylamino)-1-(4-methylphenyl)pentan-1-one [other name(s): (4- Methyl-alpha-ethylaminopentiophenone; 4-MEAP)]	7245
	(xxiv)	1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (4'-Methyl-alpha-pyrrolidinohexiophenone; MPHP)]	744 <u>6</u>
	(xxv)	1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha- Pyrrolidinoheptaphenone; PV8)]	
	(xxvi)	1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): (4'-Chloro-alpha-pyrrolidinovalerophenone; 4-chloro-alpha-PVP)]	7443
<u>SCH</u>	EDULE		
us ha (b) N	sual nam as been a arcotic l	II shall consist of the drugs and other substances, by whatever official name, common te, chemical name, or brand name designated, listed in this section. Each drug or substances the Controlled Substances Code Number set forth opposite it. Drugs: Substances, vegetable origin or chemical synthesis. Unless specifically exclusted in another schedule, any of the following substances whether produced directly the controlled substances whether the controlled substances whether produced directly the controlled substances whether the controlled substances whether produced directly the controlled substances whether the c	bstance cepted
in	directly	by extraction from substances of vegetable origin, or independently by means of chor by combination of extraction and chemical synthesis:	
(1	opiate nalbup	n and opiate, and any salt, compound, derivative, or preparation of opium or excluding apomorphine, thebaine-derived butorphanol, dextrorphan, phine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, cone and samidorphan, and their respective salts, but including the ring:	
	(i)	(1)—Raw opium	9600*
	(ii)	(2) Opium extracts	9610*
	(iii)	(3) Opium fluid extracts	_9620*
	(iv)	(4) Powdered onium	9639*

	(v)	(5)—Granulated opium	9640*
	(vi)	(6) Tincture of opium	9630*
	(vii)	(7) Codeine	9050*
	(viii)	(8) Dihydroetorphine	9334*
	(ix)	(9) Ethylmorphine	9190*
	(x)	(10) Etorphine hydrochloride	9059*
	(xi)	(11) Hydrocodone	
	(xii)	(12) Hydromorphone	9150*
	(xiii)	(13) Metopon	
	(xiv)	(14) Morphine	9300*
	(xv)	(15) Oripavine	9330*(<u>9</u> - <u>2007</u>)
	(xvi)	(16) Oxycodone	9143*
	(xvii)	(17) Oxymorphone	9652*
	(xviii)	(18) Thebaine	9333*
	(xix)	(19) Tapentadol	9780-(5- <u>20</u> 09)
	(xx)	(20) Noroxymorphone	9668 <u>-(4-2021)</u>
(2)	equiva of this	lt, compound, derivative, or preparation thereof which is chemically lent or identical with any of the substances referred to in paragraph (b) (1) section, except that these substances shall not include the isoquinoline ds of opium.*	
(3)	Opium	poppy and poppy straw.*	
<u>(4)</u>	leaves, derivat derivat	eaves (9040) and any salt, compound, derivative, or preparation of coca (including cocaine (9041) and ecgonine (9180) and their salts, isomers, ives and salts of isomers and derivatives), and any salt, compound, ive, or preparation thereof which is chemically equivalent or identical with these substances, except that the substances shall not include:	*
	<u>(i)</u>	Ddecocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine.	*
	<u>(ii)</u>	[123I]ioflupane; or	
	(i) (iii)	[18F]FP-CIT.	
	solid o	centrate or poppy straw (the crude extract of poppy straw in either liquid, r powder form which contains the phenanthrine phenanthrene alkaloids of um poppy),	9670.*

(c) **Opiates:** (Narcotic Drugs) Unless specifically excepted or unless in another schedule, any of the following opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designations:

(1) A	lfentanil	9737-(2- <u>19</u> 87)
(2) A	lphaprodine	9010*
(3) A	nileridine	9020*
(4) B	ezitramide	9800*
(5) B	ulk Dextropropoxyphene (non-dosage forms)	9273-(9- <u>19</u> 81)
(6) C	arfentanil	9743-(9- <u>19</u> 88)
	ihydrocodeine	
(8) D	iphenoxylate	9170*
(9) F	entanyl	9801*
(10)	Isomethadone	9226*
(11)	Levo-alphacetylmethadol (LAAM)	9648-(12- <u>19</u> 93)
(12)	Levomethorphan	
(13)	Levorphanol	
(14)	Metazocine	
(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	
(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- <u>19</u> 96)
(27)	Sufentanil	9740-(9- <u>19</u> 81)
(28)	Thiafentanil	9729 <u>-(4-2021</u>)
(29)	Oliceridine	9245 <u>-(5-2022</u>)
(30)	Tianeptine	<u>(5-2022)</u>

(1) A	t effect on the central nervous system:	1100			
	shetamine, its salts, optical isomers, and salts of its optical isomers				
	namphetamine, its salts, isomers, and salts of its isomers				
` /	examefetamine	`			
	metrazine and its salts nylphenidate				
(5) Meth	lylphenidate	1724			
depressa wheneve chemica	Depressants : Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:				
	barbital				
	ethimide				
` '	obarbital				
	cyclidine				
(5) Seco	barbital	2315			
f) Hallucii	nogenic Substances:				
(1) Nabi	lone	7379-(11- <u>19</u> 87			
6, or	Another Other name(s) for nabilone: (+)-trans-3-(1,1-dimethylheptyl)-6a,7,8,10,10a-hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-ne].				
U.S.	rabinol 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 <u>-(7-2019</u>			
	ate Precursor: Unless specifically excepted or unless listed in another scheo, compound, mixture, or preparation which contains any quantity of the follows:				
g) Immedi	es:				
g) Immedi material substanc	ediate precursor to Amphetamine and Methamphetamine:				
g) Immedi material substanc		- 8501-(3- <u>19</u> 80			

	(1)	1-phenylcyclohexylamine	· 7460*	
	(ii)	1-piperidinocyclohexanecarbonitrile (PCC)	8603*	
	(3) Imme	diate precursor to Fentanyl:		
	(i)	4-anilino-N-phenethylpiperidine_(ANPP)	8333*(8- <u>20</u> 10)	
	(ii)	N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl)	8366 <u>-(4-2021)</u>	
<u>SC</u>	CHEDULE	<u>; III</u>		
(a)	usual nam	III shall consist of the drugs and other substances, by whatever official name, common or me, chemical name, or brand name designated, listed in this section. Each drug or substance assigned the DEA Controlled Substances Code Number set forth opposite it.		
(b)	compound stimulant or geomet	ts: Unless specifically excepted or unless listed in another schedule, and, mixture, or preparation which contains any quantity of the following effect on the central nervous system, including its salts, isomers (whe tric), and salts of such isomers whenever the existence of such salts, is a possible within the specific chemical designation:	g substances having a ther optical, position,	
	stimul prepar Section list for	e compounds, mixtures, or preparations in dosage unit form containing ant substances listed in Schedule II which compounds, mixtures, or rations were listed on August 25, 1971, as excepted compounds under on 308.32, and any other drug of the quantitative composition shown in those drugs or which is the same except that it contains a lesser quantilled substances————————————————————————————————————	n that utity of	
		hetamine		
	(3) Chlor	phentermine	1645*	
		rmine		
	(5) Pheno	limetrazine	1615*	
(c)	compound	nts : Unless specifically excepted or unless listed in another schedule, d, mixture, or preparation which contains any quantity of the following t effect on the central nervous system:		
	(1) Any c	ompound, mixture, or preparation containing:		
	(i)	Amobarbital	2126*	
	(ii)	Secobarbital	2316*	
	(iii)	Pentobarbital	2271*	
	(iv)	Embutramide	2020*(9- <u>20</u> 06)	
		or any salt thereof and one or more other active medicinal ingredient which are not listed in any schedule.	ts	

(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.	
	Any substance which contains any quantity of a derivative of barbituric acid or ny salt thereof	2100*
(4) (Chlorhexadol	2510*
is 5	Any drug product containing gamma hydroxybutyric acid, including its salts, somers, and salts of isomers, for which an application is approved under section 05 of the Federal Food, Drug, and Cosmetic Act	
(6) K	Ketamine. its salts, isomers, and salts of isomers	- 7285-(7- <u>19</u> 99)
	Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
	ysergic acid	
	ysergic acid amide	
(9) N	Methyprylon	
(10)	<u>Sul</u> fondiethylmethane	2600*
(11)	Sulfonethylmethane	2605*
(12)	Sulfonmethane	2610*
(13)	Tiletamine and zolazepam or any salt thereof	- 7295-(3- <u>19</u> 88)
	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,-trimethylpyrazolo-[3,4-e] [1,4,]-diazepin-7(1-H)-one. flupyrazapon.	
(14)	Perampanel	2261-(11- <u>20</u> 13)
(d) Nalo	rphine	9400*
	cotic 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	
	(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	
	(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	9810*
		material, compound, mixture, or preparation containing any of the following stic drugs or their salts, as set forth below:	
	(i)	Buprenorphine(10-02 Transfer)	9064-(6- <u>19</u> 85)
	(ii)	Reserved	
(f)	compour salts, ison	Steroids: Unless specifically excepted or unless listed in another schedule, and, mixture, or preparation containing any quantity of the following substance mers, and salts of isomers whenever the existence of such salts of isomers is particular them.	s, including its possible within
	(1) Bold	enone;	<u>(9-1991)</u>
	(2) Bold	ione;	(1- <u>20</u> 10)
	(3) Chlor	rotestosterone (4-chlortestosterone);	<u>(9-1991)</u>
	(4) Clost	ebol;	<u>(9-1991)</u>
	` ′	drochlormethyltestosterone;	
		xymethyltestosterone	
	` '	drotestosterone (4-dihydrotestosterone);	
	` ′	tanolone;	
	•	estrenol;	
	(10) Fl	uoxymesterone;	·(<u>9-1991</u>)

	(11)	Formebulone (formebolone);	<u>(9-1991)</u>
	(12)	Mesterolone;	<u>(9-1991)</u>
	(13)	Methandienone, also known as Methandrostenolone;	(9-1991 <u>)</u>
	(14)	Methandranone;	<u>(9-1991)</u>
	(15)	Methandriol;	<u>(9-1991)</u>
	(16)	Methenolone;	<u>(9-1991)</u>
	(17)	Methyltestosterone;	
	(18)	Mibolerone;	
	(19)	Nandrolone;	
	(20)	19-Nor-4,9(10)-Androstadienedione	(0 1- <u>20</u> 10)
	(21)	Norethandrolone;	
	(22)	Oxandrolone;	<u>(9-1991)</u>
	(23)	Oxymesterone;	
	(24)	Oxymetholone;	
	(25)	Stanolone;	<u>(9-1991)</u>
	(26)	Stanozolol;	
	(27)	Testolactone;	<u>(9-1991)</u>
	(28)	Testosterone;	<u>(9-1991)</u>
	(29)	Trenbolone	<u>(9-1991)</u>
	(30)	Prostanozol	<u>(</u> 8- <u>20</u> 12)
	(31)	Methasterone	<u>(</u> 8- <u>20</u> 12);
		and	
	(32)	Any salt, ester, or isomer of a drug or substance described or list in this	
	pa	aragraph, if that salt, ester, or isomer promotes muscle growth	<u>(9-1991)</u>
(g)		pt anabolic steroid products: Compounds, mixtures, or preparations that cod that have been exempted by the Director Secretary :	ontain an anabolic
			NDC Number
	(1)	Andro-Estro 90-4	0536-1605
	(2)	Androgyn L.A	0456-1005
	` ′	omponent E-H in Process Pellets	
	` ′	omponent E-H in Process Granulation	· ·
	` ′	omponent TE-S in process Granulation	•
	` ′	omponent TE-S in process Pellets	•

(7) d	epANDROGYN	0456-1020
(8) D	Depo-Testadiol	0009-0253
(9) D	DEPO-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
(20)	Estratest	0032-1026
(21)	Estratest HS	
(22)	Menogen	
(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	
(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

	(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	
	(47)	Testosterone Enanthate/Estradiol Valerate injection	0182-3073
	(48)	Testosterone Enanthate/Estradiol Valerate injection	0364-6618
	(49)	Testosterone Enanthate/Estradiol Valerate injection	
	(50)	Testosterone Ophthalmic Solution	_
	(51)	Tilapia Sex Reversal Feed (investigational	Ranger, Inc
(h)	admir	inary Anabolic Steroid Implant Products: Anabolic steroid products express nistration through implants in cattle or other nonhuman species exempted by tor Secretary.	•
			NDC/DIN
		omponent E-H	
		omponent E-H	
		omponent TE-S	
		omponent T-H	
		omponent T-S	
		то	
	(7) Fi	naplix-H	12799-807-10
		naplix-S	
	(9) H	eifer-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

(40) Testoderm in process film------ Alza Corp

(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 that they be used in humans, the distribution would be subject to the criminal status despite the drugs' exempted status.		
(i) (g) Hallucinogenic substances:		
(1) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin cap a U.S. Food and Drug Administration approved drug product		
[Some other names for dronabinol: (6a R-trans)-6a,7,8, 10a-tetrahydro-trimethyl-3-pentyl-6H-dibenzo [b,d] phyran-1-ol, or (-)-delta 9-(trans)-tetrahydrocannabinol]	6, 6, 9-	
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 sched compound, mixture, or preparation containing any of the following narcotic drugs calculated as the free anhydrous base or alkaloid, in limited quantities as set for	rugs, or their salts	
(1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit		
(2) Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl propionoxybutane)		
(c) <u>Depressants</u> : Unless specifically excepted or unless listed in another schedule compound, mixture, or preparation which contains any quantity of the following including its salts, isomers, and salts of isomers whenever the existence of suc salts of isomers is possible within the specific chemical designation:	ng substances,	
(1) Alprazolam	2882-(6- <u>19</u> 82)	
(2) Barbital	2145*	
(3) Bromazepam	2748-(1- <u>19</u> 85)	
(4) Camazepam	2749-(1- <u>19</u> 85)	
(5) Chloral betaine	2460*	

(6) C	hloral hydrate	2465*
(7) C	hlordiazepoxide	2744*
(8) C	lobazam	2751-(1- <u>19</u> 85)
(9) C	lonazepam	2737*
(10)	Clorazepate	2768*
(11)	Clotiazepam	2752-(1- <u>19</u> 85)
(12)	Cloxazolam	· — /
(13)	Delorazepam	2754-(1- <u>19</u> 85)
(14)	Diazepam	
(15)	Dichloralphenazone	2467-(10- <u>20</u> 02)
(16)	Estazolam	2756-(1- <u>19</u> 85)
(17)	Ethchlorvynol	2540*
(18)	Ethinamate	2545*
(19)	Ethyl loflazepate	2758-(1- <u>19</u> 85)
(20)	Fludiazepam	2759-(1- <u>19</u> 85)
(21)	Flunitrazepam	2763-(1- <u>19</u> 85)
(22)	Flurazepam	
(23)	Fospropofol	2138-(11- <u>20</u> 09)
(24)	Halazepam	2762-(6- <u>19</u> 82)
(25)	Haloxazolam	2771-(1- <u>19</u> 85)
(26)	Ketazolam	2772-(1- <u>19</u> 85)
(27)	Loprazolam	2773-(1- <u>19</u> 85)
(28)	Lorazepam	2885*
(29)	Lormetazepam	2774-(1- <u>19</u> 85)
(30)	Mebutamate	2800*
(31)	Medazepam	2836-(1- <u>19</u> 85)
(32)	Meprobamate	2820*
(33)	Methohexital	2264*
(34)	Methylphenobarbital (mephorbarbital)	2250*
(35)	Midazolam	2884-(1- <u>19</u> 85)
(36)	Nimetazepam	2837-(1- <u>19</u> 85)
(37)	Nitrazepam	2834-(1- <u>19</u> 85)
(38)	Nordiazepam	2838-(1- 19 85)

	(39)	Oxazepam	2835*
	(40)	Oxazolam	2839*
	(41)	Paraldehyde	2585*
	(42)	Petrichloral	2591*
	(43)	Phenobarbital	2285*
	(44)	Pinazepam	2883-(1- <u>19</u> 85)
	(45)	Prazepam	
	(46)	Quazepam	
	(47)	Temazepam	
	(48)	Tetrazepam	
	(49)	Triazolam	
	(50)	Zaleplon	2781-(9- <u>19</u> 99)
	(51)	Zolpidem	
	(52)	Zopiclone	2784-(1- <u>20</u> 06)
	(53)	Alfaxalone	2731-(2- <u>20</u> 14)
	(54)	Carisoprodol	
	(55)	Tramadol	
	(56)	Suvorexant	2223-(8- <u>20</u> 14)
	(57)	Brexanolone	
	(58)	Lemborexant	
	(59)	Remimazolam	2846 <u>-(5-2022)</u>
	(60)	Daridorexant	<u> 2410</u>
(d)	follow of suc	luramine: Any material, compound, mixture, or preparation which contains a wing substances, including its salts, isomers (whether optical, position, or get ch isomers, whenever the existence of such salts, isomers, and salts of isome cenfluramine	ometric), and salts rs is possible:
(e)	(e) <u>Stimulants</u> : 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:		
	(1) C	Cathine ((+)-Norpseudeophedrine	1230-(3- <u>19</u> 88)
	(2) D	Diethylpropion	1610*

(3) Fencamfamin	1760-(3- <u>19</u> 88)
(4) Fenproporex	1575-(3- <u>19</u> 88)
(5) Lorcaserin	1625-(6-2013)
(5)(6) Mazindol	1605-(6- <u>19</u> 82)
(6)(7) Mefenorex	1580-(3- <u>19</u> 88)
(7) (<u>8)</u> Modafinil	1680-(1- <u>19</u> 99)
(8)(9) Pemoline (including organometallic complexes and chelates thereof)	1530*
(9)(10) Phentermine	1640*
(10)(11)Pipradrol	
(11)(12) Serdexmethylphenidate	<u> 1729</u>
(12)(13) (10) Sibutramine	1675-(2- <u>19</u> 98)
(14) <u>(11) Solriamfetol</u>	1650-(4-2021)
(13)(15) SPA ((-)-1-dimethylamino-1,2,diphenylethane)	1635-(9- <u>19</u> 81)
(14) (12) Lorcaserin	1625 (6-2013)
(15) (13) Solriamfetol	1650_(4-2021)

(f) <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) Pentazocine	9709-(4- <u>19</u> 79)
(2) Butorphanol	9720-(4- <u>19</u> 97)
(3) Nalbuphine	(4- <u>19</u> 97)
(4) Eluxadoline	9725-(4-2017)

SCHEDULE V

- (a) Schedule V shall consist of the drugs and other substances by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.
- (b) **Narcotic Drugs:** Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation containing any of the following narcotic drugs and their salts, as set forth below.

Reserved

(c) Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following limited quantities of narcotic drugs or salts thereof, which shall include one or more nonnarcotic active medicinal ingredients in sufficient proportion to

confer upon the compound, mixture, or preparation valuable medicinal qualities possessed by the narcotic drug alone:	es other than those
(1) Not more than 200 milligrams of codeine per 100 milliliters or per 100 gra	ms*
(2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per grams.	
(3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 1 grams.	
(4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micros of atropine sulfate per dosage unit.	
(5) Not more than 100 milligrams of opium per 100 milliliters or per 100 gram	ıs*
(6) Not more than 0.5 milligrams of difenoxin and not less than 25 microgram atropine sulfate per dosage unit.	
(d) <u>Stimulants</u> : unless specifically excepted or unless listed in another schedule, compound, mixture, or preparation which contains any quantity of the following stimulant effect on the central nervous system, including its salts, isomers and	ng substances having
(1) Pyrovalerone	1485-(3- <u>19</u> 88)
(2) Ephedrine:a -{-(Methylamino)ethyl} benzene-methanol;	(10- <u>19</u> 95)
a-{-(methylamino) ethyl}benzyl alcohol; 2-methylamino-1-phenyl-1-pro 1-phenyl-1-hydroxy-2-methylaminopropane; 1-phenyl-2- methylaminopropanol; a - hydroxy-b-methylaminopropylbenzene; a pro which occurs in the Chinese herb Ma Huang (Ephedra vulgaris, Ephedra Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra	oduct a sinica
(3) Phenylpropanolamine	
(4) Pseudoephedrine	(7- <u>20</u> 05)
Pursuant to Ark. Code Ann. § 5-64-212 as amended in 2005, this Schedule V of NOT apply to any ephedrine, phenylpropanolamine, or pseudoephedrine in liq liquid gel capsule form. However, sales limits mandated by statute shall apply ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient region.	uid, liquid capsule, or to all products with
(e) Depressants: Unless specifically exempted or excluded or unless listed in and material, compound, mixture, or preparation which contains any quantity of th having a depressant effect on the central nervous system, including its salts, is isomers:	e following substances
(1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	2782-(1- <u>20</u> 06)
(2) Lacosamide	2746-(5- <u>20</u> 09)
(3) Brivaracetam	2710 <u>-(4-2017)</u>

	(4) Lasmiditan 2790 <u>-(4-2021</u>)
	(5) Cenobamate 2720 <u>-(4-2021</u>)
	(5)(6) Ganaxolone2401
(f)	Other substances:
<u>SC</u>	CHEDULE VI**
(a)	In addition to any substance placed in Schedule VI by the <u>Director Secretary</u> 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

(a) In addition to any substance placed in Schedule VI by the <u>Director Secretary</u> 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) Mari	uana	**
(2) Tetra	hydrocannabinols, unless the tetrahydrocannabinol is:	**
(i)	(A) Contained in hemp-derived cannabidiol;	<u>(6-2020)</u>
(ii)	(B) Not more than three-tenths of one percent (0.3%) of <u>delta-9</u> tetrahydrocannabinol in the hemp-derived cannabidiol on a dry weight basis as verified by a nationally accredited laboratory for quality, purity and accuracy standards; and	<u>(6-2020)</u> ***
(iii)	(C) Not approved by the United States Food and Drug Administration for marketing as a medication;	<u>(6-2020)</u>
(3) A syı	nthetic equivalent of:	
(i)	(A) The substance contained in the Cannabis plant; or	**
(ii)	(B) The substance contained in the resinous extractives of the genus Cannabis;	<u>**</u>
class; the p manu extra salts,	a divinorum or Salvinorin A, which includes all parts of the plant presently affed botanically as Salvia divinorum, whether growing or not, the seeds of lant, any extract from any part of the plant, and every compound, afacture, derivative, mixture, or preparation of the plant, its seeds, or its cts, including salts, isomers, and salts of isomers when the existence of the isomers, and salts of isomers is possible within the specific chemical nation;	<u>**</u>

(5) Synthetic substances, derivatives, or their isomers in the chemical structural classes described below in subdivisions (a)(5)(i)-(a)(5)(x) of this section and also specific unclassified substances in subdivision (a)(5)(xi) of this section. Compounds of the structures described in this subdivision (a)(5), regardless of numerical designation of atomic positions, are included in this subdivision (a)(5).

The synthetic substances, derivatives, or their isomers included in this subdivision (a)(5) are:

<u>(i)</u>	Tetrahydrocannabinols:	
(A)	(i)Tetrahydrocannabinols, including without limitation the following:	**
8	a) (a) Delta-1 cis or trans tetrahydrocannabinol <u>[other name(s): Delta-9 cis or trans tetrahydrocannabinol]</u> , and its optical isomers;	<u>**</u>
ł	b) (b) Delta-6 cis or trans tetrahydrocannabinol <u>[other name(s): Delta-8 cis or trans tetrahydrocannabinol]</u> , and its optical isomers; and	<u>**</u>
<u>C</u>	c) (e) Delta-3.4 3.4 cis or trans tetrahydrocannabinol other name(s): Delta-6a, trans tetrahydrocannabinol, and its optical isomers—;	10a cis oı **
<u>C</u>	d) Delta-10 cis or trans tetrahydrocannabinol, and its optical isomers;	***
<u>6</u>	e) Delta-8 tetrahydrocannabinol acetate ester;	***
<u>f</u>	f)Delta-9 tetrahydrocannabinol acetate ester;	
	g) Delta-6a,10a, tetrahydrocannabinol acetate ester;	
1	h) Delta-10 tetrahydrocannabinol acetate ester; and,	***
	e)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)	(ii) Dronabinol in sesame oil and encapsulated in a soft gelatin capsule in a drug approved by the United States Food and Drug Administration is not a tetrahydrocannabinol under this subdivision (a)(5)(i);	
c c r r	(B)—Naphthoylindoles, or any compound structurally derived from 3-(1-naphtho 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:	yl)indole
(A)	(i) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**
(B)	(ii)-JWH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	**
(C)	—(iii)-JWH-018, or 1-Propyl-3-(1-naphthoyl)indole;	
(D)	(iv)-JWH-019, or 1-Hexyl-3-(1-naphthoyl)indole;	**
(E)	(v)-JWH-073, or 1-Butyl-3-(1-naphthoyl)indole;	**
(F)	(vi) JWH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
(G)	(vii) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**

(H)	(viii) JWH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl)indole; **
(I)	(ix) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole; **
(J)	(x) JWH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;*
(K)	(xi) JWH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;*
(L)	
(M	
(N)	
(O)	- (xv) EAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)-methanone; and**
(P)	(xvi) THJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3yl](naphthalen-1-yl)methanone;7024-(7-2019)
(iii)	(C)—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)	(i) JWH-175, or 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane; and**
(B)	(ii)JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1-naphthyl)methane;**
(iv)	(D)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)	(E)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)	(F) 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-

S	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)	(i)JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;*
(B)	(ii)JWH-203, or 1-Pentyl-3-(2-chlorophenylacetyl)indole;**
(C)	(iii)JWH-250, or 1-Pentyl-3-(2-methoxyphenylacetyl)indole;**
(D)	(iv)JWH-251, or 1-Pentyl-3-(2-methylphenylacetyl)indole; and**
(E)	(v)RCS-8, or 1-(2-cyclohexylethyl)-3-(2- methoxyphenylacetyl)indole; **
1 1 1	(G)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)	(i)CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]- phenol;**
(B)	(ii) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and**
(C)	(iii) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;**
s h I s	(H) ——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)	(i)AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;**
(B)	(ii) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;**
(C)	(iii)WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;**
(D)	(iv)AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole; and**
(E)	(v)—RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1H-indol-3-yl)-methanone;
I	(I) ——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)	(i)AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;
(B)	(ii) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone; **
(C)	(iii)—JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide, [other name(s): some other names: 2NE1, moved in Schedule VI in 2020]**
(D)	(iv)AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide; **
(E)	(v)5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide 7049**
(F)	(vi)STS-135, or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide;**
	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)	(i) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone; **
(B)	(ii) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl) methanone;**
(C)	(iii)A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;**
(D)	(iv)5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; **
(E)	(v)5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and**
(F)	(vi)A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]- (2,2,3,3-tetramethylcyclopropyl)methanone; **
(xi)	(K) ——Unclassified Synthetic Cannabinoids, including without limitation the following:

(A)	(i)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)	(ii)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)	(iii)—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)	(iv)—Dimethylheptylpyran or DMHP;**
(E)	(v)—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)	(vi)—URB597, or [3-(3-carbamoylphenyl)phenyl] N- Cyclohexylcarbamate; **
(G)	(vii) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one;**
(H)	(viii) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]- methanone;**
a	u) (ix)—URB602, or cyclohexyl N-(3-phenylphenyl)carbamate;**
(I)	(x)—PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3- Carboxylatecarboxylate;**
(J)	(xi)_5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;**
(K)	(xii)—BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate;**
(L)	(xiii)NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide;**
(M)	(xiv)_5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide;**
(N)	(xv)5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L-valine methyl ester 7033_(9-2018)
(O)	(xvi)MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)-L-valinate 7044-(9-2018)
(P)	(xvii)5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-carboxamido)-3,3-dimethylbutanoate; 7034-(11-2018)
(Q)	(xviii) _5-Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate 7041-(11-2018)
(R)	(xix) _MDMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3.3-dimethylbutanoate: 7042-(11-2018)

(S)	(xx)—_FUB-AMB, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate;	7021 <u>-(11-2018)</u>
(T)	(xxi) MDMB-FUBINACA, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	7020 <u>-(11-2018)</u>
(U)) (xxii)—AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-caroboxamide;	<u>7023-(7-2019)</u>
(V)) (xxiii) AB-CHMINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1 (cyclohexylmethyl)-1H-indazole-3-carboxamide;	
(W	(xxiv)(xxiv)MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide;	(11-2014)
(X)	(4-fluorobenzyl)-1H-indazole-3-carboxamide;	
(Y)) (xxvi) ADB-PINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl) 1-pentyl-1H-indazole-3-carboxamide;	
(Z)	(xxvii)5F-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	<u>7083-(6-2020)</u>
(A.	A) (xxviii) ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2 yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide	
(B)	B) (xxix)—_4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	- 7043 <u>-(4-2021)</u>
(Co	C) (xxx) 5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5 fluoropentyl)-1H-indazole-3-carboxamide	
(D)	D) (xxxi) 4-CN-CUMYL-BUTINACA, or 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7089 <u>(5-2022)</u>
(El	E) (xxxii) 5F-CUMYL-P7AICA, or 1-(5-fluoropentyl)-N-(2-phenylpropan 2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide	
(FI	F) (xxxiii) NM2201, or Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate	- 7221 <u>-(5-2022)</u>
<u>(G</u>	G) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	7036
<u>(H</u>	H) FUB-144, or (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone;	7014
<u>(II)</u>	FUB-AKB48, or N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3 carboxamide;	
A synt	thetic substance, derivative, or its isomers with:	
(i)	(A)—Similar chemical structure to any substance described in subdivision this section; or	s (a)(1)-(a)(5) of

(6) A

- (ii) (B)—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,
 - (a)(b) For persons who are twenty-one (21) years of age or older, the effective date shall be August 1, 2023.