

# **Arkansas Department of Health**

4815 West Markham Street ● Little Rock, Arkansas 72205-3867 ● Telephone (501) 661-2000 Governor Asa Hutchison Nathaniel Smith, MD, MPH, Director and State Health Officer

# SUMMARY OF PROPOSED AMENDMENTS TO RULES AND REGULATIONS 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. Items one through nine are additions made to the controlled substance list per Emergency Rule. The following items listed will include Emergency Rule additions and listed substances by current rule making procedure.

- 1. 25B-NBOH. 2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol. Page 5, (47). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (47).
- 2. 25I-NBOH. 2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol, Page 5, (48). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (48).
- 3. 5-Fluoro-ADB. methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate. Page 21, (K)(xviii). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this synthetic cannabinoid with no recognized medical use be included into Schedule VI. Page 21, (K), (xviii).

- 4. 5-Fluoro-MDMB-PICA. methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate. Page 21, (K), (xix). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this synthetic cannabinoid with no recognized medical use be included into Schedule VI. Page 21, (K), (xix).
- 5. MDMB-CHMICA. methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate. Page 21, (K)(xx). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this synthetic cannabinoid with no recognized medical use be included into Schedule VI. Page 21, (K), (xx).
- 6. FUB-AMB. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate. Page 21, (K), (xxi). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this synthetic cannabinoid with no recognized medical use be included into Schedule VI. Page 21, (K), (xxi).
- 7. MDMB-FUBINACA. methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate. Page 21, (K)(xxii). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this synthetic cannabinoid with no recognized medical use be included into Schedule VI. Page 21, (K), (xxii).
- 8. The addition of a section in Schedule V titled Other Substances page 17, (f). To incorporate the addition of a new Schedule V substance by DEA.
- 9. A drug product in finished dosage formulation that has been approved by the U.S. Food and Drug Administration that contains cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-1,3-benzenediol) derived from cannabis and no more than 0.1 percent (w/w) residual tetrahydrocannabinols. Page 17, (f), (1). To follow DEA scheduling, this drug would be included as Schedule V. Page 17, (f), (1).
- 10.AB-PINACA. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-caroboxamide. Page 21, (K)(xxiii). The DEA has scheduled this

- synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Page 21, (K), (xxiii).
- 11.AB-CHMINACA. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1- (cyclohexylmethyl)-1H-indazole-3-carboxamide. Page 21, (K), (xxiv). The DEA has scheduled this synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Page 21, (K), (xxiv).
- 12. THJ-2201. [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl) methanone. Page 18, (B), (xvi). The DEA has scheduled this synthetic cannabinoid because it has no recognized medical use. This drug would be included as Schedule VI. Page 18, (B), (xvi).
- 13.Dronabinol in an oral solution in a drug product approved for marketing by the U.S. Food and Drug Administration; [(-)-delta-9-transtetrahydrocannabinol (delta-9-THC)], Page 9, (f), (2). To follow DEA scheduling, this drug would be included as Schedule II. Page 9, (f), (2).
- 14. 25E-NBOMe. 2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine. Page 5, (49). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (49).
- 15. 25H-NBOMe. 2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine, Page 5, (50). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (50).
- 16. 25C-NBOH. 2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol. Page 5, (51). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (51).

- 17. 25H-NBOH. 2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol, Page 5, (52). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this hallucinogenic substance with no recognized medical use be included into Schedule I. Page 5, (52).
- 18. N-Ethylpentylone. 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one. Page 7, (11), (b), (18). Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this stimulant substance with no recognized medical use be included into Schedule I. Page 7, (11), (b), (18).
- 19. Two items with typos marked for clean-up.
  - Page 20, I, (ix).
  - Page 21, (K) (xvi).

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 Director 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 Director 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 Director of the Arkansas Department of Health.

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 Branch

STATE OF ARKANSAS
)
COUNTY OF SALINE
)

I, Marci Middleton, do hereby certify that Shane David, Pharm.D., well known to me, appeared before me and signed the above referenced document.

Sworn and subscribed to before me this day of , 2018.

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. Ann. \$5-64-201 and \$5-64-216.

#### SECTION II PURPOSE

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

#### SECTION III GENERAL REQUIREMENTS

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

#### SECTION IV REPEAL

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

#### CERTIFICATION

This will certify the following list of scheduling of controlled substances was adopted by the Arkansas State Board of Health at a session of the Board held in Little Rock, Arkansas on the 25th day of January, 2018, and after a Public Hearing on the 5th day of December, 2017, held in Little Rock, Arkansas, at the State Department of Health Building.

Nathaniel Smith, M.D., MPH
Director, Arkansas Department of Health

#### ARTICLE II

#### Schedule I

- (a) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) Opiates: (Narcotic Drugs) 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 paragraph (b) (34) only, the term isomer includes the optical and geometric isomers):

the o	ptical and geometric isomers):	
(1)	Acetyl-alpha-methylfentanyl(N-[1-[1-methyl-2-	
	<pre>phenethyl)-4-piperidinyl]-N-phenylacetamide)</pre>	9815-(2-86)
(2)	Acetylmethadol	9601*
(3)	Allylprodine	9602*
(4)	Alphacetylmethadol (except Levo-alphacetylmethadol	
	(LAAM)	
(5)	Alphameprodine	9604*
(6)	Alphamethadol	9605*
(7)	Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phen	yl)
	ethyl-4-piperidyl]propronanilide; 1-(1-methyl-	
	2-phenylethyl)-4(N-propanilido)piperidine)	9814-(6-82)
(8)	Alpha-methylthiofentanyl(N-[1-methyl-2-(2thienyl)	
	ethyl-4-piperidinyl]-N-phenylpropanamide)	9832-(2-86)
(9)	Benzethidine	
(10)	Betacetylmethadol	9607*
(11)	Beta-hydroxyfentanyl(N-[1-(2-hydroxy-2-phenethyl)	
	-4-piperidinyl[]-N-phenylpropanamide)	9830-(2-86)
(12)	Beta-hydroxy-3-methylfentanyl	
	[other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl	
	4-piperidinyl]-N-phenylpropamamide]	9831-(2-86)
(13)	Betameprodine	9608*
(14)	Betamethadol	
(15)	Betaprodine	
(16)	Clonitazene	
(17)	Dextromoramide	
(18)	Diampromide	9615*
(19)	Diethylthiambutene	9616*
(20)	Difenoxin	
(21)	Dimenoxadol	
(22)	Dimepheptanol	
(23)	Dimethylthiambutene	9619*
(24)	Dioxaphetyl butyrate	9621*
(25)	Dipipanone	9622*
(26)	Ethylmethylthiambutene	9623*
(27)	Etonitazene	
(28)	Etoxeridine	
(29)	Furethidine	
(30)	Hydroxypethidine	9627*
(31)	Ketobemidone	
(32)	Levomoramide	
(33)	Levophenacylmorphan	9631*
(34)	3-Methylfentanyl (N-[3-Methyl-1-(2-phenylethyl)-4-	

	piperidyl]-N-Phenylpropanamide)	9813-(10-85)
(35)	3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)	
	ethyl-4-piperidinyl]-N-phenylpropanamide)	9833-(2-86)
(36)	Morpheridine	
(37)	MPPP (1-methyl-4-phenyl-4-propionoxypiperidine)	9661-(10-85)
(38)	Noracymethadol Norlevorphanol	9633*
(39)	Normethadone	
(40)	Norpipanone	
(41) (42)	Para-fluorofentanyl (N-[4-fluorophenyl)-N-[1-(2-	9030^
(42)	phenenthyl)-4-piperindinyl]propananmide	9812-(11-86)
(43)	PEPAP 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-	9012-(11-00)
(40)	idine	9663-(10-85)
(44)	Phenadoxone	
(45)	Phenampromide	
(46)	Phenomorphan	
(47)	Phenoperidine	
(48)	Piritramide	
(49)	Proheptazine	
(50)	Properidine	9644*
(51)	Propiram	
(52)	Racemoramide	9645*
(53)	Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-	
	piperidinyl]-propanamide	9835-(2-86)
(54)	Tilidine	
(55)	Trimeperidine	9646*
(56)	Acetyl fentanyl N-(1-phenethylpiperidin-4-yl)	
	-N-phenylacetamide	-9821
(57)	Butyryl fentanyl N-(1-phenethylpiperidin-4-yl)	
	-N-phenylbutyramide	-9822
(58)	Beta-hydroxythiofentanyl N-{1-[2-hydroxy-2-	
	(thiophen-2-yl)ethyl]piperidin-4-yl}-N-	0006
(EO)	phenylpropionamide	-9836
(59)	Acetyl fentanyl 4-methylphenethyl analog N- {1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-N-	
	phenyl-acetamide monohydrochloride	
(60)	Valeryl fentanyl N-phenyl-N[1-(2-phenylethyl)	
(00)	-4-piperidinyl]-pentanamide monohydrochloride	
(61)	Furanylfentanyl N-(1-(2-phenylethyl)-	
(01)	4-piperidinyl)-N-phenylfuran-2-carboxamide	
(62)	Isobutyryl fentanyl 2-methyl-N-phenyl-N-	
(02)	[1-(2-phenylethyl)-4-piperidinyl]-propanamide	
	monohydrochloride	
(63)	Octfentanil N-(2-fluorophenyl)-2-methoxy-N-[1-(2-	
( /	phenylethyl)piperidin-4-yl]acetamide	
(64)	4-methoxy butyryl fentanyl N-(4-methoxyphenyl)-N	
	-(1-phenethylpiperidin-4-yl)butyramide monohydro	
	chloride	
(65)	Para-fluorobutyryl fentanyl N-(4-fluorophenyl)-N-	
	[1-(2-phenylethyl)-4-piperidinyl]-butanamide mono	
	hydrochloride	
(66)	Acetyl norfentanyl N-phenyl-N-4-piperidinyl-acetam	ide
	monohydrochloride	
(67)	AH-7921 3,4-dichloro-N-[(1dimethylamino)cyclohexyl	
	methyl]benzamide	
(68)	W-18 1-(4-nitrophenylethyl)piperidylidene-2-(4-	
	chlorophenyl) sulfonamide	
(69)	W-15 1-phenylethylpiperidylidene-2-(4-chlorophenyl	)
	sulfonamide	

- (70) MT-45 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine
- (71) U-47700 trans-3,4-dichloro-N-(2-(dimethylamino) cyclohexyl)-N-methylbenzamide
- (c) Opium derivatives: (Narcotic Drugs) Unless specifically excepted or unless listed in another schedule, any of the following opium derivatives, its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1)	nooccipiiino	
(2)	Acetyldihydrocodeine	9051*
(3)	Benzylmorphine	
(4)	Codeine methylbromide	9070*
(5)	Codeine-N-Oxide	9053*
(6)	Cyprenorphine	
(7)	Desomorphine	
(8)	Dihydromorphine	
(9)	Drotebanol	9335*
(10)	Etorphine (except hydrochloride salt)	9056*
(11)	Heroin	9200*
(12)	Hydromorphinol	
(13)	Methyldesorphine	
(14)	Methyldihydromorphine	9304*
(15)	Morphine methylbromide	
(16)	Morphine methylsulfonate	9306*
(17)	Morphine-N-Oxide	9307*
(18)	Myrophine	9308*
(19)	Nicocodeine	9309*
(20)	Nicomorphine	
(21)	Normorphine	
(22)	Pholcodine	
(23)	Thebacon	9315*
(24)	Mitragynine	
(25)	7-Hydroxymitragynine	

- (d) <u>Hallucinogenic substances</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation, which contains any quantity of the following hallucinogenic substance, or which contains any of its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation (for purposes of this paragraph only, the term "isomer" includes the optical, position and geometric isomers):
  - (1) alpha-ethyltryptamine ----- 7249-(12-93) 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-alpha-methylphenethylamine; 4-bromo-2,5-DMA.
  - (3) 4-bromo-2,5-dimethoxyphenethylamine ----- 7392-(8-95) 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-(3-88) Some trade or other names: DOET.
  - (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine----- 7348-(1-05) Some trade or other names: 2C-T-7.

(7)	4-methoxyamphetamineSome trade or other names: 4-methoxy-alpha-	7411*
(0)	methylphenethylamine; paramethoxyamphetamine; PMA.	7401+
(8) (9)	5-methoxy-3,4-methylenedioxy-amphetamine4-methyl-2,5-dimethoxyamphetamine	
(9)	Some trade and other names: 4-methyl-2,5-dimethoxy-alphamethylphenethylamine; "DOM"; and "STP".	7393
(10)	3,4-methylenedioxy amphetamine	7400*
(11)	3,4-methylenedioxymethamphetamine	7405-(10-85)
( ± ± /	Some trade or other names: MDMA)	, 100 (10 00)
(12)	3,4-methylenedioxy-N-ethylamphetamine	7404-(6-90)
,	Some trade or other names: N- ethy-alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA; MDE;	( )
(12)	MDEA.	7400 (6 00)
(13)	N-hydroxy-3,4-methylenedioxyamphetamine	/402-(6-90)
	Some trade or other names: N-hydroxy-alpha-methyl-3,	
(1.4.)	4 (methylenedioxy) phenethylamine; N-hydroxy MDA 3,4,5-trimethoxy amphetamine	7200*
(14)	5-methoxy-n,n-dimethyltryptamine 5-MeO-DMT	
(15) (16)	alpha-methyltryptamine	7431" (01-11)
(10)	Some trade or other names: AMT	7432 (7 03)
(17)	Bufotenine	7433*
( ± / )	Some trade and other names: 3-(beta-	7433
	Dimethylaminoethyl)-5-hydroxyindole; 3-(2-	
	dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin;	<b>:</b>
	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-(7-05)
	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	
(22)	iboga. Lysergic acid diethylamide	7215*
	Mescaline	7313**
(24)	Parahexyl	
(21)	Some trade or other names: 3-Hexyl-1-hydroxy-	7371 (7 03)
	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	<u>-</u>
	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	
(30)	Ethylamine Analog of phencyclidine	/455*
	Some trade or other names:	
	N-ethyl-1-phenylcyclohexylamine,	

```
(phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl)
     ethylamine; cyclohexamine; PCE.
     Pyrrolidine Analog of phencyclidine ----- 7458*
     Some trade or other names: 1-(1-phenylcyclohexyl)-
     pyrrolidine; PCPy; PHP
     Thiophene Analog of phencyclidine ----- 7470*
(32)
     Some trade or other names: 1-[1-(2-thienyl)
     cyclohexyl] Piperidine; 2-Thienyl analog of
     phencyclidine; TPCP; TCP.
    1-[1-(2-Thienyl)cylcohexyl]pyrrolidine ----- 7473-(9-89)
     Some other trade or other names: TCPy.
    N, N-Diallyl-5-Methoxytryptamine; 5-MeO DALT
(34)
     2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
(35)
     methyl]ethanamine; 25C-NBOMe
     2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
(36)
     methyl]ethanamine; 25I-NBOMe
(37)
    2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E---- 7509
(38)
     2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D---- 7508
     2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C--- 7519
(39)
     2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I---- 7518
(40)
     2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine;
(41)
     2C-T-2----- 7385
    2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine;
(42)
    2C-T-4----- 7532
    2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H----- 7517
(43)
     2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N---- 7521
(44)
(45)
     2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine;
     2C-P----- 7524
(46)
     2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine;
     25B-NBOMe
(47) 2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol;
     25B-NBOH
    2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol;
(48)
    25I-NBOH
    2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
    methyl]ethanamine; 25E-NBOMe
     2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)
     methyl]ethanamine; 25H-NBOMe
     2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]-
     phenol; 25C-NBOH
(52)
    2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol;
     25H-NBOH
```

- (e) <u>Depressants</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
  - (1) Phenazepam
  - (2) gamma-hydroxybutyric acid (some other names include GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate), and its known precursors and analogs -- 2010-(2-01)
  - (3) Mecloqualone ----- 2572\*
  - (4) Methagualone ----- 2565\*
  - (5) Etizolam

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

- (1) Cathinone ----- 1235-(3-88)
- (2) (<u>+</u>) CIS-4-Methylaminorex [(<u>+</u>)CIS-4,5-dihydro-4methyl-5-phenyl-2-oxazolamine] ------ 1590-(6-90)
- (3) Fenethylline ----- 1503-(9-81)
- (4) N-Benzylpiperazine----- 7493-(1-05) Some trade or other names: BZP, 1-Benzylpiperazine
- (5) N-ethylamphetamine ------ 1475-(6-82)
- (6) N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers ------ 9818-(2-86)
- (7) N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropanmide (thenylfentanyl), its optical isomers, salts and salts of isomers ------ 9834-(2-86)
- (8) N, N, Dimethylamphetamine (some other names: N, N Alpha-trimethylbenzeneethanamine; N, N, Alphatrimethylphenethylamine), its salts, optical isomers, and salts of optical isomers ----- 1480-(2-89)
- (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-(12-93)
- (10) Aminorex (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-(12-93)
- (11) Methyl-N-ethylcathinone (MEC)
- (b) Any material, compound, mixture, or preparation, whether produced directly or indirectly from a substance of vegetable origin or independently by means of chemical synthesis or by a combination of extraction and chemical synthesis, that contains any quantity of the following substances, or that contains any of the following substances' analogs, salts, isomers, and salts of isomers when the existence of the analogs, salts, isomers, and salts of isomers is possible within the specific chemical designation, with the following chemical structure is included in Schedule I:
  - (1) 4-Methylmethcathinone (Mephedrone) ----- 1248-(3-11)
  - (2) Methylenedioxypyrovalerone (MDPV) ----- (3-11)
  - (3) 3,4-Methylenedioxy-N-methylcathinone (Methylone)---- 7540-(3-11)
  - (4) 4-Methoxymethcathinone ----- (3-11)
  - (5) 3-Fluoromethcathinone ----- (3-11)
  - (6) 4-Fluoromethcathinone or ----- (3-11)
  - (7) 1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one (Butylone)
  - (8) Alpha-Pyrrolidinopentiophenone (Alpha-PVP)
  - (9) 4-methyl-N-ethylcathinone (4-MEC)
  - (10) 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)
  - (11) 2-(methylamino)-1-phenylpentan-1-one (Pentedrone)
  - (12) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone,MBDP)
  - (13) 4-fluoro-N-methylcathinone (4-FMC, Flephedrone)
  - (14) 3-fluoro-N-methylcathinone (3-FMC)
  - (15) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)
  - (16) Alpha-pyrrolidinobutiophenone ([Alpha]-PBP) or
  - (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.
- (18) N-Ethylpentylone, or 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one

#### Schedule II

- (a) Schedule II shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Controlled Substances Code Number set forth opposite it.
- (b) <u>Narcotic Drugs:</u> Substances, vegetable origin or chemical synthesis. Unless specifically excepted or unless listed in another schedule, any of the following substances whether produced directly or indirectly by extraction from substances of vegetable origin, or independently by means of chemical synthesis, or by combination of extraction and chemical synthesis:
  - (1) Opium and opiate, and any salt, compound, derivative, or preparation of opium or opiate, excluding apomorphine, dextrorphan, nalbuphine, nalmefene, naloxone and naltrexone and their respective salts, but including the following:

(1)	Raw opium	9600*
(2)	Opium extracts	9610*
(3)	Opium fluid extracts	9620*
(4)	Powdered opium	9639*
(5)	Granulated opium	9640*
(6)	Tincture of opium	9630*
(7)	Codeine	9050*
(8)	Dihydroetorphine	9334*
(9)	Ethylmorphine	
(10)	Etorphine hydrochloride	9059*
(11)	Hydrocodone	9193*
(12)	Hydromorphone	9150*
(13)	Metopon	9260*
(14)	Morphine	9300*
(15)	Oripavine	9330* (9-07)
(16)	Oxycodone	9143*
(17)	Oxymorphone	9652*
(18)	Thebaine	9333*
(19)	Tapentadol	9780-(5-09)

- (2) Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in paragraph (b) (1) of this section, except that these substances shall not include the isoquinoline alkaloids of opium.\*
- (3) Opium poppy and poppy straw.\*
- (4) Coca leaves (9040) and any salt, compound, derivative, or preparation of coca leaves, (including cocaine (9041) and ecgonine (9180) and their salts, isomers, derivatives and salts of isomers and derivatives), and any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of these substances, except that the substances shall not include decocainized

- coca leaves or extraction of coca leaves, which extractions do not
  contain cocaine or ecgonine.\*
- (5) Concentrate or poppy straw (the crude extract of poppy straw in either liquid, solid or powder form which contains the phenanthrine alkaloids of the opium 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)	Alfentanil	
(2)	Alphaprodine	9010*
(3)	Anileridine	
(4)	Bezitramide	9800*
(5)	Bulk Dextropropoxyphene (non-dosage forms)	9273-(9-81)
(6)	Carfentanil	
(7)	Dihydrocodeine	9120*
(8)	Diphenoxylate	
(9)	Fentanyl	
(10)	Isomethadone	
(11)	Levo-alphacetylmethadol (LAAM)	9648-(12-93)
(12)	Levomethorphan	9210*
(13)	Levorphanol	
(14)	Metazocine	
(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-	
	carboxylic acid	9802*
(18)	Pethidine (Meperidine)	9230*
(19)	Pethidine-Intermediate-A, 4-cyano-1-	
	methyl-4-phenylpiperidine	9232*
(20)	Pethidine-Intermediate-B, ethyl-4-	
	phenylpiperidine-4-carboxylate	9233*
(21)	Pethidine-Intermediate-C, 1-methyl-4-	
	phenylpiperidine-4-carboxylic acid	9234*
(22)	Phenazocine	9715*
(23)	Piminodine	
(24)	Racemethorphan	9732*
(25)	Racemorphan	
(26)	Remifentanil	
(27)	Sufentanil	9740-(9-81)

- (d) <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:
  - (1) Amphetamine, its salts, optical isomers, and salts of its optical isomers ------ 1100\*
     (2) Methamphetamine, its salts, isomers, and salts
  - of its isomers ------ 1105\*
  - (3) Lisdexamefetamine ----- 1205\*(7-07)
  - (4) Phenmetrazine and its salts ----- 1631\*
  - (5) Methylphenidate ----- 1724\*
- (e)  $\underline{\text{Depressants}}$ : Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any

quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Amobarbital ----- 2125\*
- Glutethimide ----- 2550-(2-91) (2)
- (3) Pentobarbital ----- 2270\*
- (4) Phencyclidine ----- 7471\*
- (5) Secobarbital ----- 2315\*

#### Hallucinogenic Substances:

- (1) Nabilone ----- 7379-(11-87) [Another name for nabilone: (+)) trans-3-(1,1dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pran-9-one].
- (2) Dronabinol in an oral solution in a drug product approved for marketing by the U.S.Food and Drug Administration; [(-)delta-9-trans-tetrahydrocannabinol(delta-9-THC)].-----7365
- Immediate Precursor: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:
  - (1)Immediate precursor to Amphetamine and Methamphetamine:
    - (i) Phenylacetone ----- 8501-(3-80) Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl Ketone; methyl benzyl Ketone.
  - Immediate precursor to Phencyclidine (PCP): (2)
    - (i) 1-phenylcyclohexylamine ----- 7460\*
    - (ii) 1-piperidinocyclohexanecarbonitrile (PCC) ----- 8603\*
  - (3) Immediate precursor to Fentyl:
    - (i) 4-anilino-n-phenethyl-4-piperidine(ANPP) ----- 8333\*(08-10)

#### Schedule III

- (a) Schedule III shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- Stimulants: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers (whether optical, position, or geometric), and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
  - (1) Those compounds, mixtures, or preparations in dosage unit form containing any stimulant substances listed in Schedule II which compounds, mixtures, or preparations were listed on August 25, 1971, as excepted compounds under Section 308.32, and any other drug of the quantitative composition shown in that list for those drugs or which is the same except that it contains a lesser quantity of controlled substances----- 1405\*

  - (2) Benzphetamine ----- 1228\*
  - (3) Chlorphentermine ----- 1645\*
  - (4) Clortermine ----- 1647\*

		1 age 10
(5)	Phendimetrazine	1615*
schedule,	epressants: Unless specifically excepted or unless list any material, compound, mixture, or preparation which cof the following substances having a depressant effect o	ontains any
nervous s	stem:	
(1)		
	(i) Amobarbital	
	(ii) Secobarbital	
	(iii) Pentobarbital (iv) Embutramide	
	or any salt thereof and one or more other	2020^(9-06)
	active medicinal ingredients which are not	
	listed in any schedule.	
(2)		
, ,	(i) Amobarbital	2126*
	(ii) Secobarbital	
	(iii) Pentobarbital	2271*
	or any salt of any of these drugs and	
	approved by the Food and Drug Administration	
(2)	for marketing only as a suppository.	
(3)	<u> </u>	2100+
(4)	derivative of barbituric acid or any salt thereof Chlorhexadol	
(5)		2310
(3)	acid, including its salts, isomers, and salts of	
	isomers, for which an application is approved under	
	section 505 of the Federal Food, Drug, and Cosmetic	
	Act	2012-(2-01)
(6)	Ketamine. its salts, isomers, and salts of isomers	7285-(7-99)
	Some other names for Ketamine: $(+-)-2-(2-$	
	Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
(7)		7300*
(8)		/310*
(9)	Methypryion	
(1)	.) Sulfonethylmethane	2605*
	2) Sulfonmethane	
	3) Tiletamine and zolazepam or any salt thereof	
	Some trade or other name for a tiletamine-	
	zolazepam combination product: Telazol.	
	Some trade or other names for tiletamine:	
	-2 (ethylamino) -2-(2-thienyl) -cyclohexanone.	
	Some trade or other names for zolazepam:	
	-4(2-fluorophenyl)-6,8-dihydro-1,3,8,-	
	<pre>trimethylpyrazolo-[3,4-e] [1,4,]-diazepin-7 (1-H)-one. flupyrazapon.</pre>	
(14	Perampanel	2261-(11-13)
	·	
(d) <u>N</u>	alorphine	9400*
(e) <u>N</u>	arcotic drugs: Unless specifically excepted or unless l	isted in another
(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	
	helow.	

(i) Not more than 1.8 grams of codeine per 100

DRAFT

below:

(10-02 Transfer)

		milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*
	(ii)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	
	(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	0007+
	(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	
	(♥)	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	
	(vi)	therapeutic amounts 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	
)	contactheir	tterial, compound, mixture, or preparation ining any of the following narcotic drugs or salts, as set forth below:	0064 (6.05)
	(1)	Buprenorphine	9064-(6-85)

(ii) Reserved

(2)

- Anabolic Steroids: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation: Items 1-28 ----- 4000-(9-91)
  - (1) Boldenone;
  - (2) Boldione; (01-10)
  - (3) Chlorotestosterone (4-chlortestosterone);
  - (4) Clostebol;
  - (5) Dehydrochlormethyltestosterone;

  - (6) Desoxymethyltestosterone (01-10)
     (7) Dihydrotestosterone (4-dihydrotestosterone);
  - (8) Drostanolone;
  - (9) Ethylestrenol;
  - (10) Fluoxymesterone;
  - (11) Formebulone (formebolone);
  - (12) Mesterolone;
  - (13) Methandienone;
  - (14) Methandranone;
  - (15) Methandriol;
  - (16) Methandrostenolone;
  - (17) Methenolone;
  - (18) Methyltestosterone;

```
(19) Mibolerone;
  (20) Nandrolone;
  (21) 19-Nor-4,9(10)-Androstadienedione (01-10)
  (22) Norethandrolone;
  (23) Oxandrolone;
  (24) Oxymesterone;
  (25) Oxymetholone;
  (26) Stanolone;
  (27) Stanozolol;
  (28) Testolactone;
  (29) Testosterone;
  (30) Trenbolone
  (31) Prostanozol-----(8-12)
  (32) Methasterone----(8-12);
  (33) Any salt, ester, or isomer of a drug or
      substance described or list in this paragraph,
      if that salt, ester, or isomer promotes muscle
      growth.
(1) Exempt anabolic steroid products: Compounds,
  mixtures, or preparations that contain an anabolic
  steroid that have been exempted by the Director:
  (1) Andro-Estro 90-4----- 0536-1605
  (2) Androgyn L.A.---- 0456-1005
  (3) Component E-H in Process Pellets----- Ivy Labs Inc.
  (4) Component E-H in Process Granulation----- Ivy Labs Inc.
  (5) Component TE-S in process Granulation----- Ivy Labs Inc.
  (6) Component TE-S in process Pellets----- Ivy Labs Inc.
     depANDROGYN----- 0456-1020
  (7)
  (8) Depo-Testadiol----- 0009-0253
  (9) 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----- 0032-1023
  (22) Menogen----- 59243-570
  (23) Menogen HS----- 59243-560
  (24) Methyltestosterone & Esterified Estrogens
      (2.5mg/1.25Mg)----- Lannett Co
  (25) Methyltestosterone & Esterified Estrogens (Half
      Strength) (1.25mg/0.625mg) ----- Lannett Co
  (26) PAN ESTRA TEST----- 0525-0175
  (27) Premarin with Methyltestosterone----- 0046-0879
  (28) Premarin with Methyltestosterone----- 0046-0878
  (29) Syntest D.S.---- 66576-231
  (30) Stntest H.S.---- 66576-230
```

		Synovex H in process bulk pellets	
	(32)	Synovex H in process granulation	
	(33)		
	(34)		
	(35)		
	(36)		
	(37)		
	(38)		
	(39)		17314-2836
	(40)		
	(41)	<u> </u>	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)		
		injection	0364-6611
	(46)	Testosterone Cypionate/Estradiol Cypionate	
		injection	0402-0257
	(47)		
		injection	0182-3073
	(48)	Testosterone Enanthate/Estradiol Valerate	
		injection	0364-6618
	(49)	Testosterone Enanthate/Estradiol Valerate	
		injection	0402-0360
	(50)	Testosterone Ophthalmic Solution	Allergan
	(51)	Tilapia Sex Reversal Feed (investigational)	Ranger, Inc.
(2)		rinary Anabolic Steroid Implant Products:	· .
		olic steroid products expressly intended for	
		nistration through implants in cattle or other	
	nonhı	uman species exempted by the Director.	
			NDC/DIN
	(1)	Component E-H	021641-002
	(2)	Component E-H	
	(3)	Component TE-S	021641-004
	(4)	Component T-H	0211641-006
	(5)	Component T-S	0211641-005
	(6)	F-TO	
	(7)	Finaplix-H	12799-807-10
	(8)	Finaplix-S	12799-807-07
	(9)	Heifer-old	Boehringer
	(10)	Heifer-old	
	(11)	Heifer-old	
	(12)	Implus-H	
	(13)	Implus-H	
	, ,	•	01968327
	(14)	Masculinizing Feed for Fish (Invesitigational) -	
	(15)	Revalor-G	
	(16)	Revalor-H	
	(17)	Revalor-S	
	(18)	Synovex H	
	(19)	Synovex H	
	(20)	Synovex Plus	
		oia Sex Reversal Feed (investigational)	
			-

If veterinary products that are granted exempted status are subsequently distributed with the intent that they be used in humans, the distribution would

be subject to the criminal sanctions of the CSA despite the drugs' exempted status.

#### (g) <u>Hallucinogenic substances</u>:

#### Schedule IV

- (a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Narcotic drugs</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:
  - (1) Not more than 1 milligram of different and not less than 25 micrograms of atropine sulfate per dosage unit.----- 9167\*
  - (2) Dextro propoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane) ----- 9278-(11-87)
- (c) <u>Depressants</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1)	Alprazolam	
(2)	Barbital	
(3)	Bromazepam	
(4)	Camazepam	
(5)	Chloral betaine	
(6)	Chloral hydrate	
(7)	Chlordiazepoxide	2744*
(8)	Clobazam	
(9)	Clonazepam	
(10)	Clorazepate	
(11)	Clotiazepam	
(12)	Cloxazolam	
(13)	Delorazepam	
(14)	Diazepam	2765*
(15)	Dichloralphenazone	
(16)	Estazolam	2756-(1-85)
(17)	Ethchlorvynol	
(18)	Ethinamate	2545*
(19)	Ethyl loflazepate	2758-(1-85)
(20)	Fludiazepam	2759-(1-85)
(21)	Flunitrazepam	
(22)	Flurazepam	
(23)	Fospropofol	2138-(11-09)
(24)	Halazepam	2762-(6-82)

(25)	Haloxazolam	2771-(1-85)
(26)	Ketazolam	
(27)	Loprazolam	
(28)	Lorazepam	
(29)	Lormetazepam	
(30)	Mebutamate	
(31)	Medazepam	
(31)	Meprobamate	
· - /	Methohexital	
(33)		
(34)	Methylphenobarbital (mephorbarbital)	
(35)	Midazolam	
(36)	Nimetazepam	
(37)	Nitrazepam	
(38)	Nordiazepam	
(39)	Oxazepam	
(40)	Oxazolam	
(41)	Paraldehyde	
(42)	Petrichloral	
(43)	Phenobarbital	
(44)	Pinazepam	
(45)	Prazepam	
(46)	Quazepam	
(47)	Temazepam	2925-(9-81)
(48)	Tetrazepam	2886-(1-85)
(49)	Triazolam	2887-(7-83)
(50)	Zaleplon	2781-(9-99)
(51)	Zolpidem	2783-(12-93)
(52)	Zopiclone	
(53)	Alfaxalone	2731-(02-14)
(54)	Carisoprodol	
(55)	Tramadol	
(56)	Suvorexant	
(00)		(0 //

- (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) Fenfluramine ----- 1670\*
- (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:
  - Cathine ((+)-Norpseudeophedrine)----- 1230-(3-88) (1)Diethylpropion ----- 1610\* (2) Fencamfamin ----- 1760-(3-88) (3) (4) Fenproporex ----- 1575-(3-88) (5) Mazindol ----- 1605-(6-82) (5) Mefenorex ----- 1580-(3-88) (6) Modafinil----- 1680-(1-99) Pemoline (including organometallic complexes and chelates thereof) ----- 1530\* Phentermine ----- 1640\* (9) (10) Pipradrol ----- 1750-(9-81) (11) Sibutramine ----- 1675-(2-98) (12) SPA ((-)-1-dimethylamino-1, 2, diphenylethane) ---- 1635-(9-81)

- (13) Lorcaserin ------ 1625-(6-13)
- (f) Other substances: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts; 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-(4-79)
  - (2) Butorphanol ----- 9720-(4-97)
  - (3) Nalbuphine ----- (4-97)
  - (4) Eluxadoline ----- 9725

#### 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.
  - \* (2) Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams.
  - \* (3) Not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 grams.
  - \* (4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit.
  - \* (5) Not more than 100 milligrams of opium per 100 milliliters or per 100 grams.
  - \* (6) Not more than 0.5 milligrams of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.
- (d) <u>Stimulants</u>: unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having stimulant effect on the central nervous system, including its salts, isomers and salts of isomers:
  - (1) Pyrovalerone ----- 1485-(3-88)
  - (2) Ephedrine:a -{-(Methylamino)ethyl}benzene-methanol; (10-95)
    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 yulgaris, Ephedra sinica Stapf.,
    - Ma Huang (Ephedra vulgaris, Ephedra sinica Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra spp.
  - (3) Phenylpropanolamine ----- (7-05)

(4) Pseudoephedrine ----- (7-05)

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

- (e) **Depressants:** Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers:

  - (2) Lacosamide ----- 2746-(05-09)
  - (3) Brivaracetam ----- 2710

#### (f) Other substances:

(1) A drug product in finished dosage formulation that has been approved by the U.S. Food and Drug Administration that contains cannabidiol(2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-1, 3-benzenediol) derived from cannabis and no more than 0.1 percent (w/w) residual tetrahydrocannabinols.

#### Schedule VI\*\*

- (a) In addition to any substance placed in Schedule VI by the Director 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
  - (3) A synthetic equivalent of:
    - (A) The substance contained in the Cannabis plant; or
    - (B) The substance contained in the resinous extractives of the genus Cannabis;
  - (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) (A)-(J) of this section and also specific unclassified substances in subdivision (a) (5) (K) 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:
- (A) (i) Tetrahydrocannabinols, including without limitation the following:
  - (a) Delta-1 cis or trans tetrahydrocannabinol, and its optical isomers;
  - (b) Delta-6 cis or trans tetrahydrocannabinol, and its optical isomers; and
  - (c) Delta-3.4 cis or trans tetrahydrocannabinol, and its optical isomers.
  - (ii) 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) (A);
- (B) Naphthoylindoles, or any compound structurally derived from 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation the following
  - (i) JWH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;
  - (ii) JWH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;
  - (iii) JWH-018, or 1-Propyl-3-(1-naphthoyl)indole;
  - (iv) JWH-019, or 1-Hexyl-3-(1-naphthoyl)indole;
  - (v) JWH-073, or 1-Butyl-3-(1-naphthoyl)indole;
  - (vi) JWH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;
  - (vii) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1naphthoyl)indole
  - (viii) JWH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl) indole;
  - (ix) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole;
  - (x) JWH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1naphthoyl)indole;
  - (xi) JWH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;
  - (xii) JWH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole;
  - (xiii) AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl) indole;
  - (xiv) MAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1-naphthalenyl)-methanone; and
  - (xv) EAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1naphthalenyl)-methanone;
  - (xvi) THJ-2201, or [1-(5-fluoropenty1)-1H-indazol-3-yl](naphthalen-1-yl)methanone;
- (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:
  - (i) JWH-175, or 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane; and
  - (ii) JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1naphthyl)methane;
- (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;
- (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;
- (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-2piperidinyl)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:
  - (i) JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;
  - (ii) JWH-203, or 1-Pentyl-3-(2-chlorophenylacetyl)indole;
  - (iii) JWH-250, or 1-Pentyl-3-(2-methoxyphenylacetyl)indole;
  - (iv) JWH-251, or 1-Pentyl-3-(2-methylphenylacetyl)indole; and
  - (v) RCS-8, or 1-(2-cyclohexylethyl)-3-(2methoxyphenylacetyl)indole;
- (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:
  - (i) CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3hydroxycyclohexyl]-phenol;
  - (ii) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and
  - (iii) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;
- (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-2piperidinyl) 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:
  - (i) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;
  - (ii) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;
  - (iii) WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;
  - (iv) AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2iodobenzoyl)indole; and

- (v) RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1Hindol-3-yl)-methanone;
- (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:
  - (i) AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;
  - (ii) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;
  - (iii) 2NE1, or 1-pentyl-3-(1-adamantylamido)indole;
  - (iv) JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide;
  - (v) AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3carboxamide;
  - (vi) 5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5fluoropentyl)-1H-indazole-3-carboxamide;
  - (vii) STS-135, or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3carboxamide;
  - (viii) MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide
  - (ix) AB-FUBINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-flurorobenzyl)-1H-indazole-3-carboxamide;
  - (x) ADB-PINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)1-pentyl-1H-indazole-3-carboxamide;
- (J) 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) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3tetramethylcyclopropyl)methanone;
  - (ii) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3tetramethylcyclopropyl)methanone;
  - (iii) A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;
  - (iv) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;
  - (v) 5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and
  - (vi) A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone; or
- (K) Unclassified Synthetic Cannabinoids, including without limitation the following:
  - (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;

- (ii) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10atetrahydrobenzo[c]chromen-1-ol;
- (iii) HU-211, or Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-6,6dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10atetrahydrobenzo[c]chromen-1-ol;
- (iv) Dimethylheptylpyran or DMHP;
- (v) WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone;
- (vii) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1benzoxazin-4-one;
- (viii)AKB-48, or N-(1-adamantyl)-1-pentylindazole-3carboxamide;
- (ix) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1-naphthalenyl]methanone;
- (x) URB602, or cyclohexyl N-(3-phenylphenyl) carbamate;
- (xi) PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3carboxylate;
- (xii) 5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;
- (xiii)BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3carboxylate;
- (xiv) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3carboxamide;
- (xv) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1Hindole-3-carboxamide;
- (xvi) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3yl]carbonyl]-L-valine methyl ester;
- (xvii) MMB\_CHMICA, or methyl-(1-cyclohexylmethyl)-1H-indole-3carbonyl)-L-valinate;
- (xviii) 5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;
- (xix) 5-Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate;
- (xx) MDMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate;
- (xxi) FUB-AMB, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate;
- (xxii) MDMB-FUBINACA, or methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;
- (xxiii) AB-PINACA. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-caroboxamide;
- (xxiv) AB-CHMINACA. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide
- (6) A synthetic substance, derivative, or its isomers with:
  - (A) Similar chemical structure to any substance described in subdivisions (a) (1)-(5) of this section; or
  - (B) Similar pharmacological effects to any substance described in subdivisions (a) (1) (5) of this section.
- (b) However, director shall not delete a controlled substance listed in this section from Schedule VI.
- (c) (1) If notice has been given to the director that the United States Food and Drug Administration has designated, rescheduled, or descheduled a marijuana-

derived substance under federal law and approved for marketing the marijuanaderived substance as a prescription medication, the director shall consider the designation, rescheduling, or descheduling of the marijuana-derived substance under this chapter.

<sup>\*-</sup>Scheduled before April, 1979.

<sup>\*\*-</sup>Schedule VI is revised to conform to Act 329 of 2013.

# QUESTIONNAIRE FOR FILING PROPOSED RULES AND REGULATIONS WITH THE ARKANSAS LEGISLATIVE COUNCIL

DEPARTMENT/AGENCY	EPARTMENT/AGENCY Arkansas Department of Health				
DIVISION	Pharmacy Services and Drug Control				
DIVISION DIRECTOR	Shane David, Pharm	.D.			
CONTACT PERSON	Shane David, Pharm	.D.			
ADDRESS	Slot 25, 4815 West N	Markham, Little R	Rock, AR 7	2205	
PHONE NO. 501 661-232 NAME OF PRESENTER AT MEETING		501 661-2769 Laura S	MAIL Shue, Gene		vid@arkansas.gov el
PRESENTER E-MAIL La	ura.Shue@arkansas.g	ov			
	INSTE	RUCTIONS			
necessary. C. If you have a method of i this Rule" below. D. Submit two (2) copies of two (2) copies of the prop  Donna K. Dav Administrativ Arkansas Leg Bureau of Leg	<ul> <li>C. If you have a method of indexing your rules, please give the proposed citation after "Short Title of this Rule" below.</li> <li>D. Submit two (2) copies of this questionnaire and financial impact statement attached to the front of two (2) copies of the proposed rule and required documents. Mail or deliver to:         <ul> <li>Donna K. Davis</li> <li>Administrative Rules Review Section</li> <li>Arkansas Legislative Council</li> </ul> </li> </ul>				
One Capitol M Little Rock, A ************************************	R 72201 ***********************************	**************************************	*****	*****	*****
2. What is the subject of the prule?	proposed	heduling of conti	rolled subs	tances.	
3. Is this rule required to comregulation?  If yes, please provide the ficitation.			Yes [		No 🔀
4. Was this rule filed under th	ne emergency provision	ons of the Admini	istrative Pr	ocedure A	.ct?
If yes, what is the effective rule?	date of the emergence	(Portion of 11/01/18—		 ffect via ei	No ⊠ mergency rule
When does the emergency expire?	rule				

	Will this emergency rule be promulgated under the permanent provision	ions of the Administrative		
	Procedure Act?	Yes 🗌	No 🖂	
5.	Is this a new rule? Yes No No If yes, please provide a brief summary explaining the regulation.	_		
	Does this repeal an existing rule? Yes No No If yes, a copy of the repealed rule is to be included with your complet replaced with a new rule, please provide a summary of the rule giving does.			
ru]	Is this an amendment to an existing le? Yes No No If yes, please attach a mark-up showing the changes in the existing rusubstantive changes. Note: The summary should explain what the up copy should be clearly labeled "mark-up."	le and a summary amendment doe	of the s, and the mark	
6.	Cite the state law that grants the authority for this proposed rule? If co Code citation. Ark. Code Ann. §§ 5-64-201 - 5-64-216	odified, please giv	e the Arkansas	
	What is the purpose of this proposed rule? Why is it necessary? odate to the List of Controlled Substances for the State of Arkansas.			
	Please provide the address where this rule is publicly accessible in electronic required by Arkansas Code § 25-19-108(b).  4p://www.healthyarkansas.com/rules_regs/rules_regs.htm	ectronic form via t	he Internet as	
9.	Will a public hearing be held on this proposed rule? Yes No If yes, please complete the following:			
	Date: 01/16/19			
	Time: 10:00 a.m.			
	Room #2512 of the Dept. of Health Bldg. at 4815 West Markham, Little Place: Rock, AR			
10	. When does the public comment period expire for permanent promulg January 16, 2019, 4:30 p.m.	•	ide a date.)	
11	. What is the proposed effective date of this proposed rule? (Must prov May 1, 2019	,		
	. Please provide a copy of the notice required under Ark. Code Ann. § blication of said notice		l proof of the	
13	. Please provide proof of filing the rule with the Secretary of State and as required pursuant to Ark. Code Ann. § 25-15-204(e).	the Arkansas Stat	e Library	

14. Please give the names of persons, groups, or organizations that you expect to comment on these rules? Please provide their position (for or against) if known. Not known

## FINANCIAL IMPACT STATEMENT

## PLEASE ANSWER ALL QUESTIONS COMPLETELY

DEPARTMENT		IMENI	Arkansas Department of Health						
DIVISION		)N	Pharmacy Services and Drug Control						
PERSON COMPLETING THIS STATEMENT Shane David, Pharm.D									
TE	TELEPHONE         501 661-2325         FAX         501 661-2769         EMAIL:         shane.david@arkansas.gov								
	To comply with Ark. Code Ann. § 25-15-204(e), please complete the following Financial Impact Statement and file two copies with the questionnaire and proposed rules.								
SE	IORT	TITLE O	F THIS RULE	List	of Controlled	d Substa	nces		
1.	Does this proposed, amended, or repealed rule have a financial impact? Yes \( \subseteq \) No \( \subseteq \)								
2.	econ	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							No 🗌	
	If an agency is proposing a more costly rule, please state the following:								
	(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 more costly rule is based on the interests of public health, safety, or welfare, and if so, please explain; and;								
	(d) Whether the reason is within the scope of the agency's statutory authority; and if so, please explain.								
4.	If the	If the purpose of this rule is to implement a federal rule or regulation, please state the following:							
	(a)	What is the	e cost to implen	nent th	e federal rule	or regu	lation?		
Current Fiscal Year				Next	Fiscal Year	<u>r</u>			
General Revenue 0 Federal Funds 0 Cash Funds 0					ral Revenue al Funds Funds	0 0			

Other (Identify)	0	Other (Identify)	0				
Total	0	Total	0				
(b) What is the	ne additional cost of the state rule?						
Current Fisca	l Year	Next Fiscal Yea	<u>r</u>				
General Revenu		General Revenue					
Federal Funds	$\frac{0}{0}$	Federal Funds	$\frac{0}{0}$				
Cash Funds	0	Cash Funds	0				
Special Revent		Special Revenue	2 0				
Other (Identify		Other (Identify)					
Other (Identity	) _0	Office (fucinity)					
Total	0	Total					
explain how the  Current Fiscal Ye  0	•	Next Fiscal Ye	<u>ear</u>				
	ontrolled and unregulated substance occur for those selling the product		eculative as to the revenue				
	implement this rule? Is this the cost of the program or grant? Please explain how the government is						
Current Fiscal Ye	ear	Next Fiscal Ye	ear				
\$ 0		\$ 0	<del></del>				
This is an unco	ntrolled and unregulated substance occur for those selling the product	e. It would be purely sp	eculative as to the revenue				
or obligation of private entity, p	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 No					
time of filing th	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;						
	the agency seeks to address with juired by statute;	the proposed rule, include	ding a statement of whether				

- (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.