# 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

#### REVISED NOTICE OF PUBLIC HEARING

The Arkansas Department of Health is extending the period to accept public comments on the Rules Pertaining to the List of Controlled Substances in Arkansas from December 22, 2023 to January 29, 2024.

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

The Arkansas Department of Health will hold a public hearing at 9:30 a.m. on January 23, 2024, at the Arkansas Department of Health, Room 2508, located at 4815 West Markham, Little Rock, Arkansas, 72205 to allow interested persons to comment on the proposed amendments.

The public may submit written comments no later than 4:30 p.m. on January 29, 2024 .Comments may be sent by email to <a href="mailto:shane.david@arkansas.gov">shane.david@arkansas.gov</a> or by mail to Shane David, Pharm.D., Pharmacy Services Section, Arkansas Department of Health, 4815 West Markham Street, Slot #25, Little Rock, Arkansas, 72205.

## 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, 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), (67), now (47) is an item that has been marked for clean-up.
- 2. Fentanyl Related Substances language has been added, marked for clean-up, and will have substances following this updated language. Schedule I, (b), (78), now (52), (i) through (vi). The language indicates the following:

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; *or*
- v. Replacement of the N-propionyl group by another acyl group;
- vi. Fentanyl-related substances shall include but are not limited to:

- 3. Schedule I, (b) (1), (7), (8), (11), (12), (34), (35), (42), (53), (56 through 65), (72 through 77) and (79 through 96) are current controlled substances. In addition, Schedule I, (b), (8) is marked for clean-up. These substances are moved to Fentanyl-related substances designation within Schedule I. Schedule I, (b), (52), (i) through (vi), (A) through (QQ). Relocation of these substances will result in subsequent numbering changes within Schedule I, (b).
- 4. Brorphine. The DEA has scheduled this opioid analgesic substance into Schedule I because it has no recognized medical use. This drug would be included in Schedule I to follow DEA. Schedule I, (b), (54).
- 5. Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested Benzimidazole-substances designation for substances with the following language be included into Schedule I without a recognized medical use.

In addition, Felisia Lackey, Chief Forensic Drug Section, Arkansas State Crime Laboratory, also requested the following benzimidazole substances also listed after implemented language which include: 4'-Hydroxy Nitazene, 5-Aminoisotonitazene, Butonitazene, Etodesnitazene [other name(s): Etazene], Flunitazene, Isotodesnitazene, Metodesnitazene, N-Desethyl Etonitazene, N-Desethyl Isotonitazene, N-Piperidinyl Etonitazene [other name(s): Etonitazepipne], N-Pyrrolidino Etonitazene, [other name(s): Etonitazepyne], N-Pyrrolidino Protonitazene and Protonitazene.

Of note, Clonitazene, Etonitazene and Isotonitazene are Schedule I substances and will be relocated from their current position and placed in the Benzimidazole-opioid substances designation within Schedule I with subsequent numbering changes.

Metonitazene. The DEA has placed this opioid analgesic substance into Schedule I because it has no recognized medical use. This drug would be in the Benzimidazole-opioid substances designation within Schedule I. The language and the following benzimidazole-opioid substances will be included into Schedule I. Schedule I, (b), (55), (i) through (vi), (A)

through (Q). The language and added substances are listed as the following:

Benzimidazole-opioid substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Benzimidazole-opioid substances means any substance not otherwise listed or excepted, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has benzimidazole ring with an ethylamine at its 1-position and benzyl group at its 2-position:

- i. With or without substitution on the benzimidazole;
- ii. With or without substitution at the ethylamine;
- *iii.* With or without inclusion of the ethylamine in a cyclic structure;
- iv. With or without substitution on the benzyl ring; or
- v. With or without replacement of the benzyl ring with an aromatic ring;
- vi. Benzimidazole-opioid substances shall include but are not limited to:
  - A. 4'-Hydroxy Nitazene;
  - B. 5-Aminoisotonitazene;
  - C. Butonitazene;
  - D. Clonitazene;
  - E. Etodesnitazene, [other name(s): Etazene];
  - F. Etonitazene;
  - G. Flunitazene;
  - H. Isotonitazene;
  - I. Isotodesnitazene
  - J. Metodesnitazene;
  - K. Metonitazene;
  - L. N-Desethyl Etonitazene;
  - M. N-Desethyl Isotonitazene;
  - N. N-Piperidinyl Etonitazene, [other name(s): Etonitazepipne];
  - O. N-Pyrrolidino Etonitazene, [other name(s): Etonitazepyne];
  - P. N-Pyrrolidino Protonitazene; and
  - Q. Protonitazene.

6. Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested this substance designation for benzodiazepine substances with the following language to be included into Schedule I without a recognized medical use. In addition, Clonazolam, Flualprazolam, Flubromazepam Flubromazolam, and Phenazepam are current Schedule 1 substances and are relocated from their current position in Schedule I and placed in the benzodiazepine-substance designation within Schedule I with subsequent numbering changes.

Bromazolam. Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this depressant substance with no recognized medical use be included into Schedule I and placed in the benzodiazepine substances designation.

Lastly, Phenazolam [other names(s): Clobromazolam]. Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested that this depressant substance with no recognized medical use be included into Schedule I and placed in the benzodiazepine substances designation.

The language and the following benzodiazepine substances is included into Schedule I. Schedule I, (e), (4), (i) through (vii). The language and added substances are noted as the following:

Benzodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Benzodiazepine substances includes any substance not otherwise listed or excepted, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has a fused 1,4-diazepine and benzene ring structure with a phenyl connected to the diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or any combination thereof. Benzodiazepine substances shall include but are not limited to:

i. Bromazolam;

- ii. Clonazolam;
- iii. Flualprazolam;
- iv. Flubromazepam;
- v. Flubromazolam;
- vi. Phenazepam; and
- vii. Phenazolam [other name(s): Clobromazolam].
- 7. Felisia Lackey, Chief Forensic Chemist-Drug Section, Arkansas State Crime Laboratory, requested this substance designation for Thienodiazepine substances with the following language included into Schedule I without a recognized medical use.

In addition, Etizolam is relocated from its current position in Schedule I and placed in the Thienodiazepine substances designation within Schedule I with subsequent numbering changes.

The language and the following Thienodiazepine substance are included into Schedule I. Schedule I, (e), (5), (i). The language and added substances are noted as the following:

Thienodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Thienodiazepine substances includes any substance not otherwise listed or excepted, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has a fused 1,4-diazepine and thiophene ring structure with a phenyl connected to the 1,4-diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or thiophene ring, any substitution(s) on the phenyl ring, or any combination thereof. Thienodiazepine substances shall include but are not limited to:

- i. Etizolam
- 8. Methiopropamine. (*N*-methyl-1-(thiophen-2-yl)propan-2-amine). The DEA has scheduled this stimulant substance into Schedule I because it has no recognized medical use. This drug would be included as Schedule I to follow DEA. Schedule I, (f), (1), (xv).

- 9. Eutylone is a Schedule I controlled substance. To follow DEA, a controlled substance code number has been set forth opposite of this substance. Schedule I, (f), (2), (xx).
- 10. Xylazine. The potential adverse health effects when abused, and increasing national prevalence of xylazine utilized as an adulterating agent to other illicit substances poses a threat to public health and safety. Although there is no current approved use in humans, xylazine is utilized in veterinarian medicine and would be included into Schedule III. Schedule III, (c), (15).
- 11. Fenfluramine is a Schedule IV substance. Schedule IV, (d), (1). To follow DEA this drug is removed from the controlled substance list with subsequent outline changes that will follow within Schedule IV.
- 12. MDMB-4en-PINACA. Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate. 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. Schedule VI, (a), (5), (xi), (JJ).
- 13. CH-PIATA. N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide. 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. Schedule VI, (a), (5), (xi), (KK).

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., Branch Chief Pharmacy Services Section
STATE OF ARKANSAS )	
COUNTY OF SALINE )	
I, Marci Middleton, do hereby certify that Shane me and signed the above referenced document.	e David, Pharm.D., well known to me, appeared before
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**

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#### **ARTICLE II**

#### **SCHEDULE I**

- (a) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Opiates: (Narcotic Drugs)</u> Unless specifically excepted or unless listed in another schedule, any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, salts is possible within the specific chemical designation (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

(1) Acetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	<del>-9815-(2</del>	
(2)(1) Acetylmethadol		
(3)(2) Allylprodine		9602*
(4)(3) Alphacetylmethadol (except Levo-alphacetylmethadol (LAAM)		9603*
(5)(4) Alphameprodine		9604*
(6)(5) 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)]	<del>- 9814-(6</del>	<del>-1982)</del>
(8) Alpha-methylthiofentanyl(N-[1-methyl-2-(2thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	<del>-9832-(2</del>	<del>-1986)</del>
(9)(6) Benzethidine		9606*
(10)(7) Betacetylmethadol		9607*
(11) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	<del>-9830-(2</del>	<del>-1986)</del>
(12) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl) 3-methyl-4-piperidinyl]-N-phenylpropamamide]		<del>-1986)</del>
(13)(8) Betameprodine		9608*
(14)(9) Betamethadol		9609*
(15)(10) Betaprodine		9611*
(16) Clonitazene		9612*
(17)(11) Dextromoramide		9613*
(18)(12) Diampromide		9615*
(19)(13) Diethylthiambutene		9616*

(20)(14) Difenoxin	9168*
(21)(15) Dimenoxadol	9617*
(22)(16) Dimepheptanol	9618*
(23)(17) Dimethylthiambutene	9619*
(24)(18) Dioxaphetyl butyrate	9621*
(25)(19) Dipipanone	9622*
(26)(20) Ethylmethylthiambutene	9623*
(27) Etonitazene	9624*
(28)(21) Etoxeridine	9625*
(29)(22) Furethidine	9626*
(30)(23) Hydroxypethidine	
(31)(24) Ketobemidone	
(32)(25) Levomoramide	9629*
(33)(26) Levophenacylmorphan	9631*
(34) 3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-piperidyl]-N-Phenylpropanamide)] 9	<del>813-(10-1985)</del>
(35) 3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	<del>9833-(2-1986)</del>
(36)(27) Morpheridine	9632*
(37)(28) MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)] 9	661-(10-1985)
(38)(29)_Noracymethadol	
(39)(30) Norlevorphanol	
(40)(31) Normethadone	9635*
	9636*
(41)(32) Norpipanone	7030
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-	<del>812 (11-1986)</del>
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-	<del>812-(11-1986)</del>
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide] (43)(33) PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine]	<del>812-(11-1986)</del> 9663-(10-
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide] (43)(33) PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] 1985)	<del>812 (11-1986)</del> 9663-(10-
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide] (43)(33) PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] 1985) (44)(34) Phenadoxone	812 (11-1986) 9663-(10- 9637* 9638*
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl) N-[1-(2-phenenthyl) 4-piperindinyl]propananmide] (43)(33) PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] 1985) (44)(34) Phenadoxone	812 (11-1986) 9663-(10- 9637* 9638* 9647*
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl) N-[1-(2-phenenthyl) 4-piperindinyl]propananmide] (43)(33) PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] 1985) (44)(34) Phenadoxone	812 (11-1986) 9663-(10 9637* 9638* 9647* 9641*

(50)(40) Properidine	9644*
(51)(41) Propiram	9649*
(52)(42) Racemoramide	9645*
(53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide	9835-(2-1986)
(54)(43) Tilidine	9750-(9-1981)
(55)(44) Trimeperidine	9646*
(56) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	<del>-9821 (4-2017)</del>
(57) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	<del>-9822-(4-2017)</del>
(58) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	<del>-9836-(4-2017)</del>
(59) Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2 (4-methylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	(4-2017)
(60) Valeryl fentanyl [other name(s): N-phenyl-N[1-(2-phenylethyl)-4-piperidinyl]-pentanamide]	<del>-9840-(4-2017)</del>
(61) Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran 2-carboxamide]	<del>-9834-(4-2017)</del>
(62) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	<del>-9827-(4-2017)</del>
(63) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	<del>-9838 (4-2017)</del>
(64) 4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin 4-yl)butyramide]	(4-2017)
(65) Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide]	<del>-9823-(4-2017)</del>
(66)(45) Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide]	(4-2017)
(67)(46) AH-7921 [other name(s): 3,4-dichloro- <i>N</i> -[(1-dimethylamino)cyclohexylmethyl]benzamide]	-9551-(4-2017)
(68)(47) W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4-chlorophenyl)sulfonamide]	(4-2017)
(69)(48) W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4-chlorophenyl)Sulfonamide]	(4-2017)
(70)(49) MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine]	9560-(4-2017)
(71)(50) U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)-N-methylbenzamide]	9547-(4-2017)

<del>(/2) Acryl ientanyl jother name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]</del>	<del>9811-(6-2020)</del>
(73) 4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin 4-yl)isobutyramide]	9824-(6-2020)
(74) Tetrahydrofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide]	9843-(6-2020)
(75) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)- N-phenylcyclopropanecarboxamide]	9845-(4-2021)
(76) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	<del>-9825-(4-2021)</del>
(77) Ortho-fluorofentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9816-(4-2021)
(78)(51) 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.	
(vi) Fentanyl-related substances shall include, but are not limited to:	
(A) Acetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	<u>-9815-(2-1986)</u>
(B) Alpha-methylfentanyl [other name(s): (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propronanilide; 1-(1-methyl-2-phenylethyl)-4(N-propanilido)piperidine)]	9814-(6-1982)
(C) Alpha-methylthiofentanyl(N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]- N-phenylpropanamide)	9832-(2-1986)
(D) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	9830-(2-1986)
(E) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide]	9831-(2-1986)

piperidyl]-N-Phenylpropanamide)]	9813-(10-1985)
(G) 3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	- 9833-(2-1986)
(H) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide]	9812-(11-1986)
(I) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamid	
(J) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9821-(4-2017)
(K) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	9822-(4-2017 <u>)</u>
(L) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	
(M) Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	(4-2017)
(N) Valeryl fentanyl [other name(s): N-phenyl-N[1-(2-phenylethyl)-4-piperidinyl]-pentanamide]	
(O) Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran-2-carboxamide]	9834-(4-2017 <u>)</u>
(P) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	<u>9827-(4-2017)</u>
(Q) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	<u>9838-(4-2017)</u>
(R) 4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	(4-2017)
(S) Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide]	9823-(4-2017)
(T) Acryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]	- <u>9811-(6-2020)</u>
(U) 4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	<u>- 9824-(6-2020)</u>
(V) Tetrahydrofuranyl [other name(s): N-(1-phenethylpiperidin-4-yl) N-phenyltetrahydrofuran-2-carboxamide]	
(W) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl) N-phenylcyclopropanecarboxamide]	
(X) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	

phenethylpiperidin-4-yl)propionamide]	9816-(4-2021)
(B)(Z) Crotonyl fentanyl [other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide]	
(C)(AA) Cyclopentyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide]	9847-(5-2022)
(D)(BB) Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9826-(5-2022)
(E)(CC) Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	9837-(5-2022)
(F)(DD) Beta-methyl fentanyl [other name(s): N-phenyl-N-(1-(2-phenylpropyl) piperidin-4-yl)propionamide]	9856-(5-2022)
(G)(EE) Beta'-phenyl fentanyl [other name: N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide]	9842-(5-2022)
(H)(FF) 2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	9855-(5-2022)
(I)(GG) 4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) piperidin-4-yl)-N-phenylacetamide]	9819-(5-2022)
(1-phenethylpiperidin-4-yl)butyramide]	9846-(5-2022)
(K)(II) Ortho-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	-9848-(5-2022)
(L)(JJ) Ortho-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9820-(5-2022)
(M)(KK)Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9817-(5-2022)
(N)(LL) Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide]	9841-(5-2022)
(O)(MM) Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide]	
(P)(NN) Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4-yl)(phenyl)carbamate]	9851-(5-2022)
(Q)(OO) Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide]	9852-(5-2022)
(R)(PP) Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9853-(5-2022)
(S)(QQ) Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide]	9854-(5-2022)

(79) Isotonitazene 9614
(80)(52) Zipeprol 9873
(81)(53) Brorphine 9098
<ul> <li>(54) Benzimidazole-opioid substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Benzimidazole-opioid substances includes any substance, not otherwise listed or excepted, and for which no exemption or approval is in effect under Section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has benzimidazole ring with an ethylamine at its 1-position and benzyl group at its 2-position: <ol> <li>With or without substitution on the benzimidazole;</li> <li>With or without substitution at the ethylamine;</li> <li>With or without inclusion of the ethylamine in a cyclic structure;</li> <li>With or without substitution on the benzyl ring; or</li> <li>With or without replacement of the benzyl ring with an aromatic ring.</li> </ol> </li> </ul>
(vi) Benzimidazole-opioid substances shall include but are not limited to:
(A)4'-Hydroxy Nitazene;
(B) 5-Aminoisotonitazene;
(C) Butonitazene;
(D) Clonitazene ;9612 *
(E) Etodesnitazene, [other name(s): Etazene];
(F) Etonitazene ; 9624*
(G) Flunitazene;
(H) Isotonitazene; 9614
(I) Isotodesnitazene;
(J) Metodesnitazene;
(K) Metonitazene; 9757
(L) N-Desethyl Etonitazene;
(M) N-Desethyl Isotonitazene;
(N) N-Piperidinyl Etonitazene [other name(s): Etonitazepipne];
(O) N-Pyrrolidino Etonitazene [other name(s): Etonitazepyne];
(P) N-Pyrrolidino Protonitazene; and
(Q) Protonitazene.

(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) A	cetorphine	9319*
(2) A	cetyldihydrocodeine	9051*
(3) B	enzylmorphine	9052*
(4) C	odeine methylbromide	9070*
· /	Sodeine-N-Oxide	
	SyprenorphineSyprenorphine	
	Desomorphine	
	Pihydromorphine	
(9) D	Protebanol	9335*
(10)	Etorphine (except hydrochloride salt)	9056*
(11)	Heroin	9200*
(12)	Hydromorphinol	9301*
(13)	Methyldesorphine	9302*
(14)	Methyldihydromorphine	
(15)	Morphine methylbromide	9305*
(16)	Morphine methylsulfonate	9306*
(17)	Morphine-N-Oxide	9307*
(18)	Myrophine	
(19)	Nicocodeine	9309*
(20)	Nicomorphine	9312*
(21)	Normorphine	9313*
(22)	Pholcodine	9314*
(23)	Thebacon	9315*
(24)	Mitragynine	(11-2015)
(25)	7-Hydroxymitragynine	(11-2015)

(d) Hallucinogenic substances: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation, which contains any quantity of the following hallucinogenic substance, or which contains any of its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation (for purposes of this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

(1) al	lpha-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	
(23)	Mescaline	7381*
(24)	Parahexyl	7374-(7-1983)
	Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl.	
(25)	Peyote	· 7415*
	Meaning all parts of the plant presently classified botanically as Lophophora williamsii Lemaire, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salts, derivative, mixture or preparation of such plant, its seeds or extracts. (Interprets 21 USC 812 (c), Schedule I (c) (12)).	
(26)	N-ethyl-3-piperidyl benzilate	7482*
(27)	N-methyl-3-piperidyl benzilate	7484*
(28)	Psilocybin	· 7437*
(29)	Psilocyn	····· 7438*
(30)	Ethylamine Analog of phencyclidine	7455*
	Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE.	
(31)	Pyrrolidine Analog of phencyclidine	7458*
	Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine: PCPy: PHP	

(32)	Thiophene Analog of phencyclidine	7470*
	Some trade or other names: 1-[1-(2-thienyl) cyclohexyl] Piperidine; 2-Thienyl analog of phencyclidine; TPCP; TCP.	
(33)	1-[1-(2-Thienyl)cylcohexyl]pyrrolidine	- 7473-(9-1989)
	Some other trade or other names: TCPy.	
(34)	N,N-Diallyl-5-Methoxytryptamine;	(6-2012)
	Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT	
(35)	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 5C-NBOMe	(5-2013)
(36)	5I-NBOMe	
(37)	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E	
(38)	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D	7508-(11-2013)
(39)	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C	7519-(11-2013)
(40)	2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I	7518-(11-2013)
(41)	2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-2	7385-(11-2013)
(42)	2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-4	7532-(11-2013)
(43)	2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H	7517-(11-2013)
(44)	2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N	7521-(11-2013)
(45)	2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine; 2C-P	7524-(11-2013)
(46) N	2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine; 25B-IBOMe	(9-2018)
(47) N	2-[[[2-(4-bromo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25B-	(11-2018)
(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH	(11-2018)
	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5E-NBOMe	(7-2019)
	2-(2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 25H-IBOMe	(7-2019)
	2-[[[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino]methyl]- phenol; 25C-	(7-2019)
(52)	2-[[[2-(2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25H-NBOH	(7-2019)
(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	- 1245-(5-2022)
	Some trade or other names: Para-methoxymethamphetamine; PMMA	
(54)	2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one	7286

Some trade or other names: Methoxetamine; MXE

(e) **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) Phenazepam	(6-2012)
(2)(1) gamma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate], and its known precursors and analogs. Precursors include but are not limited to: gamma-butyrolactone	2010 (2 2001)
(3)(2) Mecloqualone	,
(4)(3) Methaqualone	
- 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	
(5) Etizolam	``
	<del>(4-2021)</del>
(7) Flualprazolam	(4-2021)
(8) Flubromazepam	(4-2021)
(9) Flubromazolam	(4-2021)
<ul> <li>(4) Benzodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Benzodiazepine substances includes any substance, not otherwise listed or excepted, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has a fused 1,4-diazepine and benzene ring structure with a phenyl connected to the diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or any combination thereof. Benzodiazepine substances shall include but are not limited to:  (i) Bromazolam;</li> </ul>	
(ii) Clonazolam;	(4-2021)
(iii) Flualprazolam;	(4-2021)
(iv) Flubromazepam;	(4-2021)
(v) Flubromazolam;	(4-2021)
(vi) Phenazepam;	(6-2012)
(vii) Phenazolam [other name(s): Clobromazolam].	
(5) Thienodiazepine substances, their isomers, esters, ethers, salts and salts of	

isomers, esters and ethers. Thienodiazepine substances includes any substance, not otherwise listed or excepted, and for which no exemption or approval is in

effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has a fused 1,4-diazepine and thiophene ring structure with a phenyl connected to the 1,4-diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or thiophene ring, any substitution(s) on the phenyl ring, or any combination thereof. Thienodiazepine substances shall include but are not limited to:

(i) Etizolam ----- (12-2014)

## (f) Stimulants:

(1) 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	1225 (2. 1099)
salts, isomers, and salts of isomers:	1225 (2 1000)
(i) Cathinone 12	1233-(3-1988)
(ii) (±) CIS-4-Methylaminorex [(±)CIS-4,5-dihydro-4- methyl-5-phenyl-2- oxazolamine]	1590-(6-1990)
(iii) Fenethylline 1:	
(iv) N-Benzylpiperazine74	7493-(1-2005)
Some trade or other names: BZP, 1-Benzylpiperazine	
(v) N-ethylamphetamine 14	1475-(6-1982)
(vi) N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers 98	9818-(2-1986)
(vii) N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropan-mide (thenylfentanyl), its optical isomers, salts and salts of isomers 98	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	480-(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	237-(12-1993)
(x) Aminorex [other name(s): aminoraphen, 2-amino-5 phenyl-2-oxazoline, or 4,5 dihyrdo-5-phenyl-2-oxazolamine], its salts, optical isomers, and salts of optical isomers	585-(12-1993)
(xi) 4,4'-Dimethylaminorex 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	•

	(xiii)	Mesocarb	1227
	(xiv)	Methyl-N-ethylcathinone (MEC)	(6-2014)
	(xv)	Methiopropamine. [other name(s): N-methyl-1-(thiophen-2-yl)propan-2-amine]	1478
(2)	indir chem conta follo exist speci	material, compound, mixture, or preparation, whether produced directly or ectly from a substance of vegetable origin or independently by means of nical synthesis or by a combination of extraction and chemical synthesis, that aims any quantity of the following substances, or that contains any of the wing substances' analogs, salts, isomers, and salts of isomers when the ence of the analogs, salts, isomers, and salts of isomers is possible within the fice chemical designation, with the following chemical structure is included in dule I:	e
	(i)	4-Methylmethcathinone (Mephedrone)	
	(ii)	Methylenedioxypyrovalerone (MDPV)	
	(iii)	3,4-Methylenedioxy-N-methylcathinone (Methylone	
	(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)	
	(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)	,
		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)	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 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;	
	0	B) At the 3-position with an alkyl substituent; or	

	(0	C) At the nitrogen atom with alkyl or dialkyl groups, or by inclusion of the nitrogen atom in a cyclic structure.	
	(xvii	i) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one (N-Ethylpentylone) 754	3-(7-2019)
	(xix)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)754	7-(4-2021)
	(xx)	1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone ( <u>Eutylone</u> ) <u>754</u>	<del>9-</del> (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
	(xxii	i) 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)]	7446
	(xxv)	1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha-Pyrrolidinoheptaphenone; PV8)]	7548
	(xxvi	i) 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>	HEDUL	<u>e II</u>	
(a)	usual nai	e II 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 assigned the Controlled Substances Code Number set forth opposite it.	
(b)	or unless indirectly	<b>Drugs:</b> Substances, vegetable origin or chemical synthesis. Unless specifically slisted in another schedule, any of the following substances whether produced directly by extraction from substances of vegetable origin, or independently by means of s, or by combination of extraction and chemical synthesis:	ectly or
	opiat nalbı	am and opiate, and any salt, compound, derivative, or preparation of opium or the excluding apomorphine, thebaine-derived butorphanol, dextrorphan, aphine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, exone and samidorphan, and their respective salts, but including the wing:	
	(i)		9600*
	(ii)	Raw opium	
			9610*
	(iii)	Raw opium	
	(iii) (iv)	Raw opium Opium extracts	9620*
		Raw opium Opium extracts Opium fluid extracts	9620* 9639*

		(vii)	Codeine	9050*
		(viii)	Dihydroetorphine	9334*
		(ix)	Ethylmorphine	9190*
		(x)	Etorphine hydrochloride	9059*
		(xi)	Hydrocodone	9193*
		(xii)	Hydromorphone	9150*
			Metopon	
			Morphine	
			Oripavine	
			Oxycodone	
		(xvii	Oxymorphone	9652*
			i) Thebaine	
		(xix)	Tapentadol	9780-(5-2009)
		(xx)	Noroxymorphone	9668-(4-2021)
	(2)	equivof th	salt, compound, derivative, or preparation thereof which is chemically valent or identical with any of the substances referred to in paragraph (b) (1) is section, except that these substances shall not include the isoquinoline oids of opium.*	
	(3)	Opiu	am poppy and poppy straw.*	
	(4)	leave deriv deriv	a leaves (9040) and any salt, compound, derivative, or preparation of coca es, (including cocaine (9041) and ecgonine (9180) and their salts, isomers, ratives and salts of isomers and derivatives), and any salt, compound, rative, or preparation thereof which is chemically equivalent or identical with of these substances, except that the substances shall not include:	*
		(i)	Decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine;	*
		(ii)	[123I]ioflupane; or	
		(iii)	[ <sup>18</sup> F]FP-CIT.	
	(5)	solid	centrate or poppy straw (the crude extract of poppy straw in either liquid, or powder form which contains the phenanthrene alkaloids of the opium by),	9670.*
(c)	foll wh	lowin eneve	(Narcotic Drugs) Unless specifically excepted or unless in another schedulg opiates, including its isomers, esters, ethers, salts, and salts of isomers, ester the existence of such isomers, esters, ethers, and salts is possible within the designations:	s and ethers
	(1)	Alfe	ntanil	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)
(7) D	hydrocodeine	9120*
` /	Piphenoxylate	
(9) F	entanyl	
(10)	Isomethadone	
(11)	Levo-alphacetylmethadol (LAAM)	
(12)	Levomethorphan	9210*
(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	9802*
(18)	Pethidine (Meperidine)	9230*
(19)	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine	9232*
(20)	Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate	9233*
(21)	Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid	9234*
(22)	Phenazocine	9715*
(23)	Piminodine	9730*
(24)	Racemethorphan	9732*
(25)	Racemorphan	9733*
(26)	Remifentanil	- 9739-(11-1996)
(27)	Sufentanil	9740-(9-1981)
(28)	Thiafentanil	9729-(4-2021)
(29)	Oliceridine	9245-(5-2022)
(30)	Tianeptine	(5-2022)

<sup>(</sup>d) **Stimulants:** Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system:

(1) Amphetamine, its salts, optical isomers, and salts of its optical isomers	1100*
(2) Methamphetamine, its salts, isomers, and salts of its isomers	
(3) Lisdexamefetamine	1205*(7-2007)
(4) Phenmetrazine and its salts	1631*
(5) Methylphenidate	1724*
(e) <b>Depressants</b> : Unless specifically excepted or unless listed in another schedule, an compound, mixture, or preparation which contains any quantity of the following st depressant effect on the central nervous system, including its salts, isomers, and sa whenever the existence of such salts, isomers, and salts of isomers is possible with chemical designation:	abstances having a lts of isomers
(1) Amobarbital	2125*
(2) Glutethimide	2550-(2-1991)
(3) Pentobarbital	
(4) Phencyclidine	7471*
(5) Secobarbital	2315*
(f) Hallucinogenic Substances:	
(1) Nabilone	7379-(11-1987)
[Other name(s) for nabilone: (±)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10 hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one].	0a-
(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)]	
(g) <b>Immediate Precursor</b> : Unless specifically excepted or unless listed in another so material, compound, mixture, or preparation which contains any quantity of the for substances:	
(1) Immediate 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) Immediate precursor to Phencyclidine (PCP):	
(i) 1-phenylcyclohexylamine	7460*
(ii) 1-piperidinocyclohexanecarbonitrile (PCC)	8603*
(3) Immediate precursor to Fentanyl:	
(i) 4-anilino-N-phenethylpiperidine (ANPP)	8333*(8-2010)

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

#### **SCHEDULE III**

- (a) Schedule III shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) **Stimulants:** Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers (whether optical, position, or geometric), and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
- (c) **Depressants**: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system:
  - (1) Any compound, mixture, or preparation containing:

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

(2) Any suppository dosage form containing:

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

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

	` /	any substance which contains any quantity of a derivative of barbituric acid or my salt thereof	2100*
		hlorhexadol	
	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	Letamine. its salts, isomers, and salts of isomers	- 7285-(7-1999)
		Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
		ysergic acid	
	(8) L	ysergic acid amide	7310*
	(9) N	Methyprylon	
	(10)	Sulfondiethylmethane	
	(11)	Sulfonethylmethane	
	(12)	Sulfonmethane	2610*
	(13)	Tiletamine and zolazepam or any salt thereof	- 7295-(3-1988)
		Some trade or other name for a tiletamine- zolazepam combination product: Telazol.	
		Some trade or other names for tiletamine: -2(ethylamino)-2-(2-thienyl)-cyclohexanone.	
		Some trade or other names for zolazepam: -4(2-fluorophenyl)-6,8-dihydro-1,3,8,-trimethylpyrazolo-[3,4-e] [1,4,]-diazepin-7(1-H)-one. flupyrazapon.	
	(14)	Perampanel	2261-(11-2013)
	<u>(15)</u>	Xylazine	
(d)	) Nalo	rphine	9400*
(e)	Nar	cotic drugs: Unless specifically excepted or unless listed in another schedule:	
	n	any material, compound, mixture, or preparation containing any of the following arcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in mited quantities as set forth below:	
	<b>(</b> i	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*

	(ii)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9804*
	(iii)	Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9807*
	(iv)	Not more than 300 milligrams of ethylmorphine per 100 milliliters or not more than 15 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9808*
	(v)	Not more than 500 milligrams of opium per 100 milliliters or per 100 grams or not more than 25 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9809*
	(vi)	Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9810*
		material, compound, mixture, or preparation containing any of the following otic drugs or their salts, as set forth below:	
	(i)	Buprenorphine 90 (10-02 Transfer)	064-(6-1985)
	(ii)	Reserved	
	` ′	Tess, via	
(f)	Anaboli compou salts, iso the spec	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:	including its ssible within 4000
(f)	Anaboli compour salts, iso the spece (1) Bold	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:	including its ssible within 4000 (9-1991)
(f)	Anaboli compour salts, iso the spece (1) Bold	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:	including its ssible within 4000 (9-1991)
(f)	Anaboli compous salts, iso the spece (1) Bolo (2) Bolo (2)	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:	including its saible within 4000 (9-1991) (1-2010)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positific chemical designation:  denone;  denone;	including its ssible within 4000 (9-1991) (1-2010) (9-1991)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo (4) Clos	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);	including its ssible within 4000 (9-1991) (1-2010) (9-1991) (9-1991)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);  stebol;	including its saible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh (6) Deso	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);  stebol;  ydrochlormethyltestosterone;	including its saible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (1-2010)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh (6) Dese (7) Dihy	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is possific chemical designation:  denone;  dione;  orotestosterone (4-chlortestosterone);  ydrochlormethyltestosterone;  oxymethyltestosterone	including its sible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
(f)	Anabolic compoursalts, iso the spect (1) Bold (2) Bold (3) Chld (4) Clos (5) Deh (6) Desc (7) Dihy (8) Dros	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);  ydrochlormethyltestosterone;  oxymethyltestosterone  ydrotestosterone (4-dihydrotestosterone);	including its sible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
(f)	Anabolic compoursalts, iso the spect (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh (6) Deso (7) Dihy (8) Dros (9) Ethy	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);  ydrochlormethyltestosterone;  oxymethyltestosterone	including its sible within 4000 (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991) (9-1991)
(f)	Anaboli compour salts, iso the spece (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh (6) Deso (7) Dihy (8) Dros (9) Ethy (10) F	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is positic chemical designation:  denone;  dione;  orotestosterone (4-chlortestosterone);  stebol;  oxymethyltestosterone  ydrotestosterone (4-dihydrotestosterone);  stanolone;  vlestrenol;	including its sible 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)
(f)	Anaboli compour salts, iso the spect (1) Bolo (2) Bolo (3) Chlo (4) Clos (5) Deh (6) Deso (7) Dihy (8) Dros (9) Ethy (10) F (11) F	c Steroids: Unless specifically excepted or unless listed in another schedule, any nd, mixture, or preparation containing any quantity of the following substances, omers, and salts of isomers whenever the existence of such salts of isomers is posific chemical designation:  denone;  denone;  orotestosterone (4-chlortestosterone);  stebol;  cydrochlormethyltestosterone;  oxymethyltestosterone  ydrotestosterone (4-dihydrotestosterone);  stanolone;  luoxymesterone;  luoxymesterone;	including its ssible 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)

	(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	(1-2010)
	(21)	Norethandrolone;	` /
	(22)	Oxandrolone;	
	(23)	Oxymesterone;	(9-1991)
	(24)	Oxymetholone;	(9-1991)
	(25)	Stanolone;	` ,
	(26)	Stanozolol;	
	(27)	Testolactone;	
	(28)	Testosterone;	(9-1991)
	(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 ragraph, if that salt, ester, or isomer promotes muscle growth	(9-1991)
(g)		pt anabolic steroid products: Compounds, mixtures, or preparations that cod that have been exempted by the Secretary:	ntain an anabolic
		NDC Number	
	(1) A	ndro-Estro 90-4	0536-1605
	(2) A	ndrogyn L.A	0456-1005
	(3) C	omponent E-H in Process Pellets	Ivy Labs Inc.
	(4) C	omponent E-H in Process Granulation	Ivy Labs Inc
	(5) Co	omponent TE-S in process Granulation	Ivy Labs Inc
	` /	omponent TE-S in process Pellets	•
		pANDROGYN	•
	(8) D	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	2) Testosterone Cypionate/Estradiol Cypionate injection	54274-530
(43	3) Testosterone Cypionate/Estradiol Cypionate injection	0182-3069
(44	Testosterone Cyp 50 Estradiol Cyp 2	0814-7737
(45	5) Testosterone Cypionate/Estradiol Cypionate injection	0364-6611
(46	Testosterone Cypionate/Estradiol Cypionate injection	0402-0257
(47	7) Testosterone Enanthate/Estradiol Valerate injection	0182-3073
(48		
(49	7) Testosterone Enanthate/Estradiol Valerate injection	0402-0360
(50	,	
(51	Tilapia Sex Reversal Feed (investigational	Ranger, Inc
` /	terinary Anabolic Steroid Implant Products: Anabolic steroid products expre ministration through implants in cattle or other nonhuman species exempted by NDC/DIN	•
(1)	Component E-H	021641-002
(2)		
(3)		
\ /	Component T-H	
(5)		0211641-005
( )	F-TO	
` ′	Finaplix-H	
` ′	Finaplix-S	
` '	Heifer-old	
(10		Č
(11		<u> </u>
(12		•
(13	•	
`	01968327	
(14	Masculinizing Feed for Fish (Invesitigational)	Rangen,Inc.
(15		_
(16		
(17	<i>'</i>	
(18	<i>'</i>	
, ,	<b>,</b>	<del>-</del>

(2	20)	Synovex Plus	0856-3904
(2	21)	Tilapia Sex Reversal Feed (investigational)	Zeigier Bros.
tl	hat tl	erinary products that are granted exempted status are subsequently distributed they be used in humans, the distribution would be subject to the criminal sanction te the drugs' exempted status.	
(i) <u>F</u>	Hallu	<u>icinogenic substances</u> :	
(		ronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in U.S. Food and Drug Administration approved drug product	
		[Some other names for dronabinol: (6a R-trans)-6a,7,8, 10a-tetrahydro-6, 6, 9 trimethyl-3-pentyl-6H-dibenzo [b,d] phyran-1-ol, or (-)-delta 9-(trans)-tetrahydrocannabinol]	-
<u>SCH</u>	IEDI	ULE IV	
u	ısual	dule IV shall consist of the drugs and other substances, by whatever official name, chemical name or brand name designated, listed in this section. Each deen assigned the DEA Controlled Substances Code Number set forth opposite in	rug or substance
c	omp	otic drugs: Unless specifically excepted or unless listed in another schedule, a ound, mixture, or preparation containing any of the following narcotic drugs, clated as the free anhydrous base or alkaloid, in limited quantities as set forth be	or their salts
(		ot more than 1 milligram of difenoxin and not less than 25 micrograms of ropine sulfate per dosage unit	9167*
(2	2) D pr	extro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-ropionoxybutane)	- 9278-(11-1987)
c ii	omp nclud	essants: Unless specifically excepted or unless listed in another schedule, any ound, mixture, or preparation which contains any quantity of the following subding its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation:	ostances,
(	1) A	lprazolam	2882-(6-1982)
(2	2) B	arbital	2145*
(.	3) B	romazepam	2748-(1-1985)
(4	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	2887-(7-1983)
(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	2400-(4-2021)
(58)	Lemborexant	2245-(4-2021)
(59)	Remimazolam	2846-(5-2022)
(60)	Daridorexant	2410
follov	uramine: Any material, compound, mixture, or preparation which ving substances, including its salts, isomers (whether optical, position is somers, whenever the existence of such salts, isomers, and salts	on, or geometric), and salts
<del>(1) F</del>	enfluramine	1670*
stimu or ge	Stimulants: Unless specifically excepted or unless listed in anoth ound, mixture, or preparation which contains any quantity of the foliant effect on the central nervous system, including its salts, isomerometric), and salts of such isomers whenever the existence of such sers is possible within the specific chemical designation:	ellowing substances having a se (whether optical, position,
(1) C	athine ((+)-Norpseudeophedrine	1230-(3-1988)
	iethylpropion	
(3) F	encamfamin	1760-(3-1988)
(4) F	enproporex	1575-(3-1988)

(5) L	orcaserin	1625-(6-2013)
(6) M	fazindol	1605-(6-1982)
(7) N	lefenorex	1580-(3-1988)
(8) M	Iodafinil	1680-(1-1999)
(9) P	emoline (including organometallic complexes and chelates thereof)	1530*
(10)	Phentermine	1640*
(11)	Pipradrol	1750-(9-1981)
(12)	Serdexmethylphenidate	1729
(13)	Sibutramine	1675-(2-1998)
(14)	Solriamfetol	
(15)	SPA ((-)-1-dimethylamino-1,2,diphenylethane)	1635-(9-1981)
subst	Other substances: Unless specifically excepted or unless listed in anotherial, compound, mixture, or preparation which contains any quantity of the ances, including its salts; isomers whether optical, position, or geometric), ers, whenever the existence of such salts, isomers, and salts of isomers is p	following and salts of such
	entazocine	
(2) B	utorphanol	9720-(4-1997)
	albuphine	
(4) E	luxadoline	9725-(4-2017)
<b>SCHED</b>	ULE 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:	ıg	
(1) Pyrovalerone 1485-(3-19	988)	
(2) Ephedrine:a -{-(Methylamino)ethyl}benzene-methanol;(10-19	995)	
a-{-(methylamino) ethyl} benzyl alcohol; 2-methylamino-1-phenyl-1-propanol; 1-phenyl-1-hydroxy-2-methylaminopropane; 1-phenyl-2-methylaminopropanol; a - hydroxy-b-methylaminopropylbenzene; a product which occurs in the Chinese herb Ma Huang (Ephedra vulgaris, Ephedra sinica Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra spp.		
(3) Phenylpropanolamine(7-20		
(3) Phenylpropanolamine(7-20) (4) Pseudoephedrine		
	005) , or th	
(4) Pseudoephedrine (7-20 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, liquid gel capsule form. However, sales limits mandated by statute shall apply to all products wit ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless of the dosa	oos), or the age	
<ul> <li>(4) Pseudoephedrine</li></ul>	one of the age	
<ul> <li>(4) Pseudoephedrine</li></ul>	one of the large o	
<ul> <li>(4) Pseudoephedrine</li></ul>	005)  or th age  onces  006) 009)	
(4) Pseudoephedrine	005)  , or th age  nces  006) 009) 017)	
(4) Pseudoephedrine	005) , or th age nces 006) 009) 017)	
(4) Pseudoephedrine	005) , or th age  006) 009) 017) 021)	

# (f) Other substances:

(1) None.

# SCHEDULE VI \*\*\*\*

(a)	under indired by a consubstant	r § 5 ectly com- anc iso	on to any substance placed in Schedule VI by the Secretary of the Department of 5-64-214, any material, compound, mixture, or preparation, whether produced by from a substance of vegetable origin or independently by means of chemical abination of extraction and chemical synthesis, that contains any quantity of the es, or that contains any of their salts, isomers, and salts of isomers when the extenses, and salts of isomers is possible within the specific chemical designation, et VI:	directly or synthesis, or e following istence of the
	(1) M	1ari	juana	**
			hydrocannabinols, unless the tetrahydrocannabinol is:	
	(i		Contained in hemp-derived cannabidiol;	
	(i	i)	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	-(6-2020) ***
	(i	ii)	Not approved by the United States Food and Drug Administration for marketing as a medication;	(6-2020)
	(3) A	syı	nthetic equivalent of:	
	(i	(i)	The substance contained in the Cannabis plant; or	**
	(i	i)	The substance contained in the resinous extractives of the genus Cannabis;	**
	cl th m ex	lass ne p nanu xtra alts,	if a divinorum or Salvinorin A, which includes all parts of the plant presently ified 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 gnation;	**
	cl sp C m T	lass peci com ume the s	hetic substances, derivatives, or their isomers in the chemical structural es described below in subdivisions (a)(5)(i)-(a)(5)(x) of this section and also affic unclassified substances in subdivision (a)(5)(xi) of this section. pounds of the structures described in this subdivision (a)(5), regardless of erical designation of atomic positions, are included in this subdivision (a)(5). synthetic substances, derivatives, or their isomers included in this subdivision are:  Tetrahydrocannabinols:	

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

	a)	Delta-1 cis or trans tetrahydrocannabinol [other name(s): Delta-9 cis or trans tetrahydrocannabinol], and its optical isomers;	**
	b)	Delta-6 cis or trans tetrahydrocannabinol [other name(s): Delta-8 cis or trans tetrahydrocannabinol], and its optical isomers;	**
	c)	Delta- 3,4 cis or trans tetrahydrocannabinol [other name(s): Delta-6a,10a cis or trans tetrahydrocannabinol], and its optical isomers;	**
	d)	Delta-10 cis or trans tetrahydrocannabinol, and its optical isomers;	***
	e)	Delta-8 tetrahydrocannabinol acetate ester;	***
	f)	Delta-9 tetrahydrocannabinol acetate ester;	***
	g)	Delta-6a,10a, tetrahydrocannabinol acetate ester;	
	h)	Delta-10 tetrahydrocannabinol acetate ester; and,	***
	i)	A product derived from industrial hemp that was produced as a result of a synthetic chemical process that converted the industrial hemp or a substance contained in industrial hemp into Delta-8, Delta-9, Delta 6a,10a, or Delta-10 tetrahydrocannabinol including their respective	
		acetate esters	***
	dru Ad	onabinol in sesame oil and encapsulated in a soft gelatin capsule in a ag product approved by the United States Food and Drug lministration is not a tetrahydrocannabinol under this subdivision (5)(i);	**
(ii)	nap the cyc (4- ind	othoylindoles, or any compound structurally derived from 3-(1-phthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at a nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-morpholinyl)ethyl group, whether or not further substituted in the lole ring to any extent and whether or not substituted in the naphthyl g to any extent, including without limitation the following:	**
	(A)JW	7H-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**
	(B) JW	7H-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	**
	(C) JW	7H-018, or 1-Propyl-3-(1-naphthoyl)indole;	
	(D) JW	H-019, or 1-Hexyl-3-(1-naphthoyl)indole;	**
	(E) JW	/H-073, or 1-Butyl-3-(1-naphthoyl)indole;	**
	(F) JW	H-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(G) JW	H-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**
	(H)JW	H-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl)indole;	**
	(I) JW	H-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole;	**
	(J) JW	7H-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;	**

	(K) JWH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl) indole;	**
	(L) JWH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole;	**
	(M) AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole;	
	(N) MAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1-naphthalenyl)-methanone;	
	(O) EAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)-methanone; and	**
	(P) THJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl)methanone;7	024-(7-2019)
(iii)	Naphthylmethylindoles, or any compound structurally derived from an H-indol-3-yl-(1-naphthyl) methane by substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation the following:	**
	(A) JWH-175, or 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane; and	**
	(B) JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1-naphthyl)methane;	**
(iv)	Naphthoylpyrroles, or any compound structurally derived from 3-(1-naphthoyl)pyrrole by substitution at the nitrogen atom of the pyrrole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the pyrrole ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-307, or (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-ylmethanone;	**
(v)	Naphthylmethylindenes, or any compound structurally derived from 1-(1-napthylmethyl)indene with substitution at the 3-position of the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-176, or E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane;	**
(vi)	Phenylacetylindoles, or any compound structurally derived from 3-phenylacetylindole by substitution at the nitrogen atom of the indole ring with alkyl, haloalkyl, alkenyl, cycloalkylmethyl,cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
	(A) JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;	**

(B) JWH-203, or 1-Pentyl-3-(2-chlorophenylacetyl)indole;	**
(C) JWH-250, or 1-Pentyl-3-(2-methoxyphenylacetyl)indole;	**
(D) JWH-251, or 1-Pentyl-3-(2-methylphenylacetyl) indole; and	
(E) RCS-8, or 1-(2-cyclohexylethyl)-3-(2- methoxyphenylacetyl)indole;	
(vii) Cyclohexylphenols, or any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not substituted in the cyclohexyl ring to any extent, including without limitation the following:	**
(A) CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol;	**
(B) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and	**
(C) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;	**
(viii) Benzoylindoles, or any compound structurally derived from a 3- (benzoyl)indole structure with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
(A) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;	**
(B) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;	**
(C) WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;	**
(D) AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole; and	**
(E) RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1H-indol-3-yl)-methanone;	**
(ix) Adamantoylindoles, or Adamantoylindazoles, including Adamantyl Carboxamide Indoles and Adamantyl Carboxamide Indazoles, or any compound structurally derived from 3-(1-adamantoyl) indole, 3-(1-adamantoyl) indazole, or 3-(2-adamantoyl)indole by substitution at a nitrogen atom of the indole or indazole ring with alkyl, haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl,cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole or indazole ring to any extent and whether or not	

	limitation the following:	**
(	(A) AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;	**
(	(B) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;	**
(	(C) JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide [other name(s): 2NE1, moved in Schedule VI in 2020]	**
(	(D) AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide;	**
(	(E) 5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	- 7049**
(	(F) STS-135, or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide;	**
(x)	Tetramethylcyclopropylcarbonylindoles or any compound structurally derived from 3-(2,2,3,3-tetramethylcyclopropylcarbonyl) indole by substitution at the nitrogen atom of the indole ring with alkyl,haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl, cycloalkylethyl, (N-methylpiperidin-2-yl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole ring to any extent, including without limitation the following:	**
(	(A) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(B) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	
(	(C) A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(D) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
(	(E) 5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and	**
(	(F) A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
(xi)	Unclassified Synthetic Cannabinoids, including without limitation the following:	**
(	(A) CP 50556-1 hydrochloride, or [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthridin-1-yl] Acetate;	**
(	(B) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;	**

(C) HU-211, or Dexanabinol,(6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;**
(D) Dimethylheptylpyran or DMHP;**
(E) WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone;**
(F) URB597, or [3-(3-carbamoylphenyl)phenyl] N-Cyclohexylcarbamate;**
(G) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one;**
(H) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]-methanone;**
a) URB602, or cyclohexyl N-(3-phenylphenyl)carbamate;**
(I) PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3-carboxylate;**
(J) 5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;**
(K) BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate;**
(L) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide; **
(M) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide;**
(N) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L-valine methyl ester 7033-(9-2018)
(O) MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)- L-valinate 7044-(9-2018)
(P) 5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-carboxamido)-3,3-dimethylbutanoate; 7034-(11-2018)
(Q) 5-Fluoro-MDMB-PICA, or methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate7041-(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) MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)- 1-(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;	(9-2018)
	(Z) 5F-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7083-(6-2020)
	(AA) ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide	7010-(4-2021)
	(BB) 4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	7043-(4-2021)
	(CC) 5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	7025-(5-2022)
	(DD) 4-CN-CUMYL-BUTINACA, or 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7089 –(5-2022)
	(EE) 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)
	(GG) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	7036
	(HH) FUB-144, or (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone;	7014
	(II) FUB-AKB48, or N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide;	7047
	(JJ)MDMB-4en-PINACA, or Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate;	
	(KK) CH-PIATA, or N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide.	
	(6) A synthetic substance, derivative, or its isomers with:	
	(i) Similar chemical structure to any substance described in subdivisions (a)(1)-(a)(5) of this section; or	**
	(ii) Similar pharmacological effects to any substance described in subdivisions (a)(1)-(a)(5) of this section	**
(b)	However, except as provided under subsection (c) of this section, the Secretary shall necontrolled substance listed in this section from Schedule VI.	ot delete a
(c)	A prescription drug approved by the United States Food and Drug Administration und 355 is excluded from Schedule VI unless the secretary objects under § 5-64-201.	er 21 U.S.C. §

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

- \*\*-Schedule VI is revised to conform to Act 329 of 2013.
- \*\*\* Schedule VI is revised further to conform to Act 629 of 2023. Each substance added to the Controlled Substances List pursuant to Act 629 of 2023 shall have the following effective dates:
  - (a) For persons who are under twenty-one (21) years of age, the effective date shall be the effective date of Act 629 of 2023; and,
  - (b) For persons who are twenty-one (21) years of age or older, the effective date shall be August 1, 2023.

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



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., Branch Chief Pharmacy Services Section
STATE OF ARKANSAS )	
COUNTY OF SALINE )	
I, Marci Middleton, do hereby certify that Shane Dame and signed the above referenced document.	avid, Pharm.D., well known to me, appeared before
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**

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#### **ARTICLE II**

# **SCHEDULE I**

- (a) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Opiates: (Narcotic Drugs)</u> Unless specifically excepted or unless listed in another schedule, any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, salts is possible within the specific chemical designation (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

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

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

	of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that is turally related to fentanyl by one or more of the following modifications:	(4-2021)
(i)	Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;	
(ii)	Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iii)	Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;	
(iv)	Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; or	
(v)	Replacement of the N-propionyl group by another acyl group.	
(vi)	Fentanyl-related substances shall include, but are not limited to:	
(.	A) Acetyl-alpha-methylfentanyl [other name(s): (N-[1-[1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide)]	9815-(2-1986)
(-	B) Alpha-methylfentanyl [other name(s): (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propronanilide; 1-(1-methyl-2-phenylethyl)-4(N-propanilido)piperidine)]	- 9814-(6-1982)
(	C) Alpha-methylthiofentanyl(N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	- 9832-(2-1986)
(.	D) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	- 9830-(2-1986)
(.	E) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide]	- 9831-(2-1986)
	F) 3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4-piperidyl]-N-Phenylpropanamide)]	9813-(10-1985)
(	G) 3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	- 9833-(2-1986)
(	H) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-4-piperindinyl]propananmide]	9812-(11-1986)
(	I) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamid	
(.	J) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9821-(4-2017)
(.	K) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	
(	L) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	

(M) Acetyl fentanyl 4-methylphenethyl analog [other name(s): N-{1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-N-phenyl-acetamide]	(4-2017)
(N) Valeryl fentanyl [other name(s): N-phenyl-N[1-(2-phenylethyl)-4-piperidinyl]-pentanamide]	-9840-(4-2017)
(O) Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran-2-carboxamide]	-9834-(4-2017)
(P) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	-9827-(4-2017)
(Q) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	-9838-(4-2017)
(R) 4-methoxy butyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	(4-2017)
(S) Para-fluorobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide]	-9823-(4-2017)
(T) Acryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide]	9811-(6-2020)
(U) 4-Fluoroisobutyryl fentanyl [other name(s): N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9824-(6-2020)
(V) Tetrahydrofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide]	9843-(6-2020)
(W) Cyclopropyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide]	9845-(4-2021)
(X) Methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide]	9825-(4-2021)
(Y) Ortho-fluorofentanyl [other name(s): N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide]	9816-(4-2021)
(Z) Crotonyl fentanyl [other name(s): (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide]	9844-(5-2022)
(AA) Cyclopentyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide]	9847-(5-2022)
(BB) Para-chloroisobutyryl fentanyl [other name(s): N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide]	9826-(5-2022)
(CC) Para-methoxybutyryl fentanyl [other name(s): N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide]	9837-(5-2022)
(DD) Beta-methyl fentanyl [other name(s): N-phenyl-N-(1-(2-phenylpropyl) piperidin-4-yl)propionamide]	9856-(5-2022)
(EE) Beta'-phenyl fentanyl [other name: N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide]	

	(FF) fluo	2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2-rophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide]	9855-(5-2022)
	(GG) pipe	4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) ridin-4-yl)-N-phenylacetamide]	9819-(5-2022)
	(HH) (1-p	Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N-henethylpiperidin-4-yl)butyramide]	9846-(5-2022)
		no-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1-methylpiperidin-4-yl)acetamide]	-9848-(5-2022)
		no-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2-nylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide]	9820-(5-2022)
	(KK) pher	Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1-methylpiperidin-4-yl)propionamide]	9817-(5-2022)
	(LL)	Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-nylbenzamide]	
	(MM) N-pl	Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-nenylthiophene-2-carboxamide]	9839-(5-2022)
	(NN) yl)(p	Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4-phenyl)carbamate]	9851-(5-2022)
	(OO) pher	Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1-nethylpiperidin-4-yl)acrylamide]	9852-(5-2022)
	(PP) N-(1	Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-phenethylpiperidin-4-yl)isobutyramide]	9853-(5-2022)
	(QQ) (1-p	Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N-henethylpiperidin-4-yl)furan-2-carboxamide]	9854-(5-2022)
(52)			
(53)	Brorphin	e	9098
ise su ar A	omers, est obstance, r oproval is ct [21 U.S	dazole-opioid substances, their isomers, esters, ethers, salts and salts of ers and ethers. Benzimidazole-opioid substances includes any not otherwise listed or excepted, and for which no exemption or in effect under Section 505 of the Federal Food, Drug, and Cosmetic J.C. 355], that structurally has benzimidazole ring with an ethylamine tion and benzyl group at its 2-position:	f
(i)	) With o	or without substitution on the benzimidazole;	
(ii	i) With o	or without substitution at the ethylamine;	
(ii	ii) With o	or without inclusion of the ethylamine in a cyclic structure;	
(ir	v) With o	or without substitution on the benzyl ring; or	
(v	) With o	or without replacement of the benzyl ring with an aromatic ring.	
(v	i) Benzi	midazole-opioid substances shall include but are not limited to:	

	(A)4'-Hydroxy Nitazene;	
	(B) 5-Aminoisotonitazene;	
	(C) Butonitazene;	
	(D) <u>Clonitazene</u> ;	9612 *
	(E) Etodesnitazene, [other name(s): Etazene];	
	(F) Etonitazene;	9624*
	(G) Flunitazene;	
	(H) Isotonitazene;	9614
	(I) Isotodesnitazene;	
	(J) Metodesnitazene;	
	(K) Metonitazene;	9757
	(L) N-Desethyl Etonitazene ;	
	(M) N-Desethyl Isotonitazene;	
	(N) N-Piperidinyl Etonitazene [other name(s): Etonitazepipne];	
	(O) N-Pyrrolidino Etonitazene [other name(s): Etonitazepyne];	
	(P) N-Pyrrolidino Protonitazene; and	
	(Q) Protonitazene.	
schee the e	um derivatives: (Narcotic Drugs) Unless specifically excepted or unless listed dule, any of the following opium derivatives, its salts, isomers, and salts of isomers is possible within the specificantion:	ners whenever
(1) A	Acetorphine	9319*
(2) A	Acetyldihydrocodeine	9051*
(3) E	Benzylmorphine	9052*
(4) (	Codeine methylbromide	9070*
(5)	Codeine-N-Oxide	9053*
(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	9302*
(14)	Methyldihydromorphine	9304*
(15)	Morphine methylbromide	9305*
(16)	Morphine methylsulfonate	9306*
(17)	Morphine-N-Oxide	9307*
(18)	Myrophine	9308*
(19)	Nicocodeine	
(20)	Nicomorphine	
(21)	Normorphine	
(22)	Pholcodine	
(23)	Thebacon	
(24)	Mitragynine	, , ,
(25)	7-Hydroxymitragynine	(11-2015)
	etric isomers):  pha-ethyltryptamine	7249-(12-1993)
(1) [[	Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-	(12 1993)
(2) 4-	ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET. bromo-2,5-dimethoxy-amphetamine	7301*
(2) न-	Some trade or other names: 4-bromo-2,5-dimethoxy-alphamethylphenethylamine; 4-bromo-2,5-DMA.	7371
(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		
(0) 2,	5-dimethoxy-4-(n)-propylthiophenethylamine	7348-(1-2005)

(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	
(16)	alpha-methyltryptamine	- 7432-(7-2005)
	Some trade or other names: AMT	
(17)	Bufotenine	7433*
	Some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine.	
(18)	Diethyltryptamine	· 7434*
	Some trade or other names: N,N-Diethyltryptamine;DET.	
(19)	Dimethyltryptamine	····· 7435*
	Some trade or other names: DMT	
(20)	5-methoxy-N,N-diisopropyltryptamine	- 7439-(7-2005)
	Some trade or other names: 5-MeO-DIPT.	
(21)	Ibogaine	· 7260*
	Some trade and other names: 7-Ethyl-6,6 beta; 7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido [1',2': 1,2] azepino [5,4-b] indole; Tabernanthe iboga.	
(22)	Lysergic acid diethylamide	7315*

(23)	Mescaline	7381*
(24)	Parahexyl	- 7374-(7-1983)
	Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl.	
(25)	Peyote	7415 <b>*</b>
	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	
(27)	N-methyl-3-piperidyl benzilate	
(28)	Psilocybin	
(29)	Psilocyn	
(30)	Ethylamine Analog of phencyclidine	7455*
	Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE.	
(31)	Pyrrolidine Analog of phencyclidine	7458*
	Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCPy; PHP	
(32)	Thiophene Analog of phencyclidine	7470*
	Some trade or other names: 1-[1-(2-thienyl) cyclohexyl] Piperidine; 2-Thienyl analog of phencyclidine; TPCP; TCP.	
(33)	1-[1-(2-Thienyl)cylcohexyl]pyrrolidine	- 7473-(9-1989)
	Some other trade or other names: TCPy.	
(34)	N,N-Diallyl-5-Methoxytryptamine;	(6-2012)
	Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT	
(35)	2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 5C-NBOMe	
(36)	2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5I-NBOMe	(8-2013)
(37)	2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine; 2C-E	7509-(11-2013)
(38)	2-(2,5-Dimethoxy-4-methylphenyl)ethanamine; 2C-D	7508-(11-2013)
(39)	2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine; 2C-C	7519-(11-2013)
(40)	2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine; 2C-I	7518-(11-2013)
(41)	2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-2	7385-(11-2013)

	(42)	2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine; 2C-T-4	7532-(11-2013)
	(43)	2-(2,5-Dimethoxyphenyl)ethanamine; 2C-H	7517-(11-2013)
	(44)	2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine; 2C-N	7521-(11-2013)
	(45)	2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine; 2C-P	7524-(11-2013)
	(46) N	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-BOH	(11-2018)
	(48)	2-[[[2-(4-iodo-2,5-dimethoxyphenyl)ethyl]amino]methyl]-phenol; 25I-NBOH-	(11-2018)
	(49) 25	2-(4-ethyl-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl) methyl]ethanamine; 5E-NBOMe	(7-2019)
	(50) N	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	(7-2019)
	(53)	1-(4-methoxyphenyl)-N-methylpropan-2-amine	1245-(5-2022)
		Some trade or other names: Para-methoxymethamphetamine; PMMA	
	(54)	2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one	7286
		Some trade or other names: Methoxetamine; MXE	
(e)	comp depre when	essants: Unless specifically excepted or unless listed in another schedule, any nound, mixture, or preparation which contains any quantity of the following subsessant 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 the ical designation:	tances having a of isomers
	hy ar	amma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4-ydroxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate], nd its known precursors and analogs. Precursors include but are not limited to:	-2010-(2-2001)
	(2) N	Mecloqualone	2572*
	(3) M	lethaqualone	2565*
	es of	enzodiazepine substances, their isomers, esters, ethers, salts and salts of isomers sters and ethers. Benzodiazepine substances includes any substance, not therwise listed or excepted, and for which no exemption or approval is in effect ander section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355],	,

on the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or any combination thereof. Benzodiazepine substances shall include but are not limited to:

(i)	Bromazolam;		

(ii)	Clonazolam;		(4-2021)	)
------	-------------	--	----------	---

- (iii) Flualprazolam:-----(4-2021)
- (iv) Flubromazepam; -----(4-2021)
- (v) Flubromazolam; -----(4-2021)
- (vi) Phenazepam; -----(6-2012)
- (vii) Phenazolam [other name(s): Clobromazolam].
- (5) Thienodiazepine substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers. Thienodiazepine substances includes any substance, not otherwise listed or excepted, and for which no exemption or approval is in effect under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], that structurally has a fused 1,4-diazepine and thiophene ring structure with a phenyl connected to the 1,4-diazepine ring, with any substitution(s) or replacement(s) on the 1,4-diazepine or thiophene ring, any substitution(s) on the phenyl ring, or any combination thereof. Thienodiazepine substances shall include but are not limited to:
  - (i) Etizolam ----- (12-2014)

#### (f) Stimulants:

(1) 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:

(i)	Cathinone	 	 1235-(	(3-1988)
(1)	Cammone	 	1233-(	(3-1766)

- (ii) (<u>+</u>) CIS-4-Methylaminorex [(<u>+</u>)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, Nemethylcathinone, methycathinone, AL-464, AL-422, AL-463 and UR-1431), its salts, optical isomers and salts of optical isomers	
	(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	1585-(12-1993)
	,	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)
		Amineptine	
	(xiii)	Mesocarb	1227
	(xiv)	Methyl-N-ethylcathinone (MEC)	(6-2014)
	(xv)	Methiopropamine. [other name(s): N-methyl-1-(thiophen-2-yl)propan-2-amine]	1478
(2)	indire chem conta follow existe speci	material, compound, mixture, or preparation, whether produced directly or ectly from a substance of vegetable origin or independently by means of call synthesis or by a combination of extraction and chemical synthesis, that any quantity of the following substances, or that contains any of the wing substances' analogs, salts, isomers, and salts of isomers when the ence of the analogs, salts, isomers, and salts of isomers is possible within the fic chemical designation, with the following chemical structure is included in dule I:	
	(i)	4-Methylmethcathinone (Mephedrone)	1248-(3-2011)
	(ii)	Methylenedioxypyrovalerone (MDPV)	(3-2011)
	(iii)	3,4-Methylenedioxy-N-methylcathinone (Methylone	7540-(3-2011)
	(iv)	4-Methoxymethcathinone	(3-2011)
	(v)	3-Fluoromethcathinone	(3-2011)
	(vi)	4-Fluoromethcathinone	(3-2011)
	(vii)	1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone)	7541-(11-2014)
	(viii)	Alpha-Pyrrolidinopentiophenone (Alpha-PVP)	7545-(11-2015)
	(ix)	4-methyl-N-ethylcathinone (4-MEC)	1249-(9-2018)
	(x)	4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP)	7498-(9-2018)
	(xi)	2-(methylamino)-1-phenylpentan-1-one (Pentedrone)	1246-(9-2018)

(XII) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk-MBDP)	7542-(9-2018)
(xiii) 4-fluoro-N-methylcathinone (4-FMC, Flephedrone)	,
(xiv) 3-fluoro-N-methylcathinone (3-FMC)	
(xv) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone)	1258-(9-2018)
(xvi) Alpha-pyrrolidinobutiophenone ([Alpha]-PBP)	
(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	e) 7543-(7-2019)
(:)1(121 1: 15 1)2(11 1: )	
(xix) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)	7547-(4-2021)
(xix) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone)	
	7549-(4-2021)
(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone ( <u>Eutylone</u> )(xxi) 2-(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone;	7549-(4-2021) 7246
<ul> <li>(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (<u>Eutylone</u>)</li></ul>	7549-(4-2021) 7246 7544
<ul> <li>(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (<u>Eutylone</u>)</li></ul>	7549-(4-2021) 7246 7544 7245
<ul> <li>(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (<u>Eutylone</u>)</li></ul>	7549-(4-2021) 7546 7544 7245
<ul> <li>(xx) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone)</li></ul>	7549-(4-2021)7246754472457548

## **SCHEDULE II**

- (a) Schedule II shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Controlled Substances Code Number set forth opposite it.
- (b) Narcotic Drugs: Substances, vegetable origin or chemical synthesis. Unless specifically excepted or unless listed in another schedule, any of the following substances whether produced directly or

indirectly by extraction from substances of vegetable origin, or independently by means of chemical synthesis, or by combination of extraction and chemical synthesis:

(1) Opium and opiate, and any salt, compound, derivative, or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, naltrexone and samidorphan, and their respective salts, but including the following:

(i) Raw opium	
(ii) Opium extracts	
(iii) Opium fluid extracts	
(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	9300*
(xv) Oripavine	
(xvi) Oxycodone	
(xvii)Oxymorphone	9652*
(xviii) Thebaine	9333*
(xix) Tapentadol	9780-(5-2009)
(xx) Noroxymorphone	9668-(4-2021)

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

a:	erivative, or preparation thereof which is chemically equivalent or identically ny of these substances, except that the substances shall not include:	W1th *
(i		is do
(i	i) [123] lioflupane; or	
(i	ii) [ <sup>18</sup> F]FP–CIT.	
S	Concentrate or poppy straw (the crude extract of poppy straw in either liquid, olid or powder form which contains the phenanthrene alkaloids of the opium oppy),	
followher	tes: (Narcotic Drugs) Unless specifically excepted or unless in another soluting opiates, including its isomers, esters, ethers, salts, and salts of isomers, ever the existence of such isomers, esters, ethers, and salts is possible within itical designations:	esters and ethers
(1) A	.lfentanil	9737-(2-1987)
(2) A	alphaprodine	9010*
(3) A	nileridine	9020*
(4) B	sezitramide	9800*
(5) B	sulk Dextropropoxyphene (non-dosage forms)	9273-(9-1981)
(6) C	arfentanil	9743-(9-1988)
(7) D	Pihydrocodeine	9120*
	oiphenoxylate	
(9) F	entanyl	9801*
(10)	Isomethadone	9226*
(11)	Levo-alphacetylmethadol (LAAM)	9648-(12-1993)
(12)	Levomethorphan	9210*
(13)	Levorphanol	9220*
(14)	Metazocine	9240*
(15)	Methadone	9250*
(16)	Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane	9254*
(17)	Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-	
C	arboxylic acid	9802*
(18)	Pethidine (Meperidine)	9230*
(19)	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine	9232*
(20)	Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate	9233*

	(21)	Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid	9234*
	(22)	Phenazocine	9715*
	(23)	Piminodine	9730*
	(24)	Racemethorphan	9732*
	(25)	Racemorphan	9733*
	(26)	Remifentanil	9739-(11-1996)
	(27)	Sufentanil	
	(28)	Thiafentanil	
	(29)	Oliceridine	
	(30)	Tianeptine	(5-2022)
(a)	comp	<b>clants:</b> Unless specifically excepted or unless listed in another schedule, around, mixture, or preparation which contains any quantity of the following lant effect on the central nervous system:  Imphetamine, its salts, optical isomers, and salts of its optical isomers	substances having a
		ethamphetamine, its salts, isomers, and salts of its isomers	
		sdexamefetamine	
	(4) Pł	nenmetrazine and its salts	1631*
	(5) M	ethylphenidate	1724*
(e)	depres when	essants: Unless specifically excepted or unless listed in another schedule, abound, mixture, or preparation which contains any quantity of the following sant effect on the central nervous system, including its salts, isomers, and sever the existence of such salts, isomers, and salts of isomers is possible wincal designation:	substances having a salts of isomers
	(1) A	mobarbital	2125*
	(2) G	lutethimide	2550-(2-1991)
	(3) Pe	entobarbital	2270*
	(4) Pl	nencyclidine	7471*
	(5) Se	ecobarbital	2315*
(f)	Hallu	cinogenic Substances:	
	(1) N	abilone	7379-(11-1987)
		[Other name(s) for nabilone: (±)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10, hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one].	10a-

	U.S.	nabinol in an oral solution in a drug product approved for marketing by the Food and Drug Administration; [(-)-delta-9-trans-thydrocannabinol(delta-9-THC)]	7365-(7-2019)
(g)		iate Precursor: Unless specifically excepted or unless listed in another scheol, compound, mixture, or preparation which contains any quantity of the followers:	<u> </u>
	(1) Imm	nediate 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	nediate precursor to Phencyclidine (PCP):	
	(i)	1-phenylcyclohexylamine	7460*
	(ii)	1-piperidinocyclohexanecarbonitrile (PCC)	8603*
	(3) Imn	nediate 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>	HEDUL		
(a)	usual na	te III shall consist of the drugs and other substances, by whatever official name, chemical name, or brand name designated, listed in this section. Each draw assigned the DEA Controlled Substances Code Number set forth opposite it	rug or substance
(b)	compou stimular or geom	ants: Unless specifically excepted or unless listed in another schedule, any mend, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system, including its salts, isomers (whether operic), and salts of such isomers whenever the existence of such salts, isomers is possible within the specific chemical designation:	tances having a ptical, position,
	stim prep Sect list	ose compounds, mixtures, or preparations in dosage unit form containing any sulant substances listed in Schedule II which compounds, mixtures, or parations were listed on August 25, 1971, as excepted compounds under aion 308.32, and any other drug of the quantitative composition shown in that for those drugs or which is the same except that it contains a lesser quantity of crolled substances————————————————————————————————————	
	(2) Ben	zphetamine	1228*
	(3) Chlo	orphentermine	1645*
	(4) Clo	termine	1647*
	(5) Phen	ndimetrazine	1615*

aepres	ssant effect on the central nervous system:	
(1) Ar	ny compound, mixture, or preparation containing:	
(i)	Amobarbital	212
(ii)	) Secobarbital	231
(iii		
(iv	y) Embutramide	2020*(9-200
	or any salt thereof and one or more other active medicinal ingredients which are not listed in any schedule.	
(2) Ar	ny suppository dosage form containing:	
(i)	Amobarbital	212
(ii)		231
(iii	i) Pentobarbital	227
	or any salt of any of these drugs and approved by the Food and Drug Administration for marketing only as a suppository.	
	ny substance which contains any quantity of a derivative of barbituric acid or y salt thereof	210
(4) Ch	nlorhexadol	251
isc	ny drug product containing gamma hydroxybutyric acid, including its salts, omers, and salts of isomers, for which an application is approved under section of the Federal Food, Drug, and Cosmetic Act	- 2012-(2-20
(6) Ke	etamine. its salts, isomers, and salts of isomers	7285-(7-19
	Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
(7) Ly	vsergic acid	730
(8) Ly	sergic acid amide	<b></b> 731
(9) Me	ethyprylon	257
(10)	Sulfondiethylmethane	260
(11)	Sulfonethylmethane	260
(12)	Sulfonmethane	261
(13)	Tiletamine and zolazepam or any salt thereof	- 7295-(3-19
	Some trade or other name for a tiletamine- zolazepam combination product: Telazol.	

			ome trade or other names for tiletamine: -2(ethylamino)-2-(2-thienyl)-yclohexanone.	
			ome trade or other names for zolazepam: -4(2-fluorophenyl)-6,8-dihydro-,3,8,-trimethylpyrazolo-[3,4-e] [1,4,]-diazepin-7(1-H)-one. flupyrazapon.	
	(14	) P	Perampanel	2261-(11-2013)
	(15	5) X	Xylazine	
(d)	Na	lorph	ine	9400*
(e)	Na	arcot	ic drugs: Unless specifically excepted or unless listed in another schedule:	
	(1)	narc	material, compound, mixture, or preparation containing any of the following totic drugs, or their salts calculated as the free anhydrous base or alkaloid, in ted quantities as set forth below:	
		(i)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*
		(ii)	Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9804*
		(iii)	Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9807*
		(iv)		
		(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	
		(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*
	(2)	•	material, compound, mixture, or preparation containing any of the following totic drugs or their salts, as set forth below:	
		(i)	Buprenorphine(10-02 Transfer)	- 9064-(6-1985)
		(ii)	Reserved	
(f)	cor salt	npou ts, iso	ic Steroids: Unless specifically excepted or unless listed in another schedule, and, mixture, or preparation containing any quantity of the following substance omers, and salts of isomers whenever the existence of such salts of isomers is sific chemical designation:	es, including its possible within

(1) Boldenone;	(9-1991)
(2) Boldione;	(1-2010)
(3) Chlorotestosterone (4-chlortestosterone);	(9-1991)
(4) Clostebol;	(9-1991)
(5) Dehydrochlormethyltestosterone;	(9-1991)
(6) Desoxymethyltestosterone	(1-2010)
(7) Dihydrotestosterone (4-dihydrotestosterone);	
(8) Drostanolone;	(9-1991)
(9) Ethylestrenol;	(9-1991)
(10) Fluoxymesterone;	(9-1991)
(11) Formebulone (formebolone);	(9-1991)
(12) Mesterolone;	(9-1991)
(13) Methandienone, also known as Methandrostenolone;	(9-1991)
(14) Methandranone;	(9-1991)
(15) Methandriol;	
(16) Methenolone;	(9-1991)
(17) Methyltestosterone;	
(18) Mibolerone;	(9-1991)
(19) Nandrolone;	(9-1991)
(20) 19-Nor-4,9(10)-Androstadienedione	
(21) Norethandrolone;	
(22) Oxandrolone;	(9-1991)
(23) Oxymesterone;	(9-1991)
(24) Oxymetholone;	(9-1991)
(25) Stanolone;	
(26) Stanozolol;	(9-1991)
(27) Testolactone;	(9-1991)
(28) Testosterone;	(9-1991)
(29) Trenbolone	(9-1991)
(30) Prostanozol	(8-2012)
(31) Methasterone	(8-2012);
and	

(32) 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	(9-1991)			
(g) Exempt anabolic steroid products: Compounds, mixtures, or preparations that contain an anabolic steroid that have been exempted by the Secretary:				
NDC Number				
(1) Andro-Estro 90-4	0536-1605			
(2) Androgyn L.A				
(3) Component E-H in Process Pellets	Ivy Labs Inc.			
(4) Component E-H in Process Granulation	-			
(5) Component TE-S in process Granulation	_			
(6) Component TE-S in process Pellets	Ivy Labs Inc			
(7) depANDROGYN	0456-1020			
(8) Depo-Testadiol				
(9) DEPO-T.E				
(10) depTESTROGEN				
(11) Duomone	52047-360			
(12) DUO-SPAN II				
(13) DURATESTRIN	43797-016			
(14) Essian				
(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			

Premarin with Methyltestosterone ------ 0046-0878

(28)

(29)	Syntest D.S	66576-231
(30)	Stntest H.S	66576-230
(31)	Synovex H in process bulk pellets	Syntex Animal
(32)	Synovex H in process granulation	Syntex Animal
(33)	Synovex Plus in process granulation	Fort Dodge
(34)	Synovex Plus in process bulk pellets	Fort Dodge
(35)	TEST-ESTRO Cypionates	
(36)	Testoderm with Adhesive 4mg/d	
(37)	Testoderm 4mg/d	
(38)	Testoderm 6mg/d	17314-4609
(39)	Testoderm with Adhesive 6mg/d	
(40)	Testoderm in process film	Alza Corp
(41)	Testoderm with Adhesive in process film	Alza Corp
(42)	Testosterone Cypionate/Estradiol Cypionate injection	54274-530
(43)	Testosterone Cypionate/Estradiol Cypionate injection	0182-3069
(44)	Testosterone Cyp 50 Estradiol Cyp 2	0814-7737
(45)	Testosterone Cypionate/Estradiol Cypionate injection	0364-6611
(46)	Testosterone Cypionate/Estradiol Cypionate injection	0402-0257
(47)	Testosterone Enanthate/Estradiol Valerate injection	0182-3073
(48)	Testosterone Enanthate/Estradiol Valerate injection	0364-6618
(49)	Testosterone Enanthate/Estradiol Valerate injection	0402-0360
(50)	Testosterone Ophthalmic Solution	Allergan
(51)	Tilapia Sex Reversal Feed (investigational	Ranger, Inc
	rinary Anabolic Steroid Implant Products: Anabolic steroid products expinistration through implants in cattle or other nonhuman species exempted NDC/DIN	-
(1) (	Component E-H	021641-002
(2) (	Component E-H	01968327
(3) (	Component TE-S	021641-004
(4) (	Component T-H	0211641-006
(5) (	Component T-S	0211641-005
(6) I	G-TO	00093351

(7) Fi	naplix-H	12799-807-10
(8) Fi	naplix-S	12799-807-07
(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.
<ul><li>(14)</li><li>(15)</li></ul>	Revalor-G	12799-811
` /		12799-811
(15)	Revalor-G	12799-811 12799-810
(15) (16)	Revalor-G	12799-811 12799-810 12799-809 0856-3901
(15) (16) (17)	Revalor-G	12799-811 12799-810 12799-809 0856-3901
(15) (16) (17) (18)	Revalor-G	12799-811 12799-810 12799-809 0856-3901 Syntex

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

### (i) Hallucinogenic substances:

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

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

### **SCHEDULE IV**

- (a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Narcotic drugs</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:

(1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit 9167*
(2) Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-propionoxybutane) 9278-(11-1987)
(c) <u>Depressants</u> : Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
(1) Alprazolam 2882-(6-1982)
(2) Barbital 2145*
(3) Bromazepam 2748-(1-1985)
(4) Camazepam 2749-(1-1985)
(5) Chloral betaine2460*
(6) Chloral hydrate 2465*
(7) Chlordiazepoxide 2744*
(8) Clobazam 2751-(1-1985)
(9) Clonazepam 2737*
(10) Clorazepate 2768*
(11) Clotiazepam 2752-(1-1985)
(12) Cloxazolam 2753-(1-1985)
(13) Delorazepam 2754-(1-1985)
(14) Diazepam 2765*
(15) Dichloralphenazone 2467-(10-2002)
(16) Estazolam 2756-(1-1985)
(17) Ethchlorvynol 2540*
(18) Ethinamate 2545*
(19) Ethyl loflazepate 2758-(1-1985)
(20) Fludiazepam 2759-(1-1985)
(21) Flunitrazepam 2763-(1-1985)
(22) Flurazepam 2767*
(23) Fospropofol 2138-(11-2009)
(24) Halazepam 2762-(6-1982)
(25) Haloxazolam 2771-(1-1985)

(26)	Ketazolam	2772-(1-1985)
(27)	Loprazolam	2773-(1-1985)
(28)	Lorazepam	2885*
(29)	Lormetazepam	2774-(1-1985)
(30)	Mebutamate	2800*
(31)	Medazepam	2836-(1-1985)
(32)	Meprobamate	2820*
(33)	Methohexital	2264*
(34)	Methylphenobarbital (mephorbarbital)	2250*
(35)	Midazolam	2884-(1-1985)
(36)	Nimetazepam	2837-(1-1985)
(37)	Nitrazepam	2834-(1-1985)
(38)	Nordiazepam	
(39)	Oxazepam	2835*
(40)	Oxazolam	
(41)	Paraldehyde	2585*
(42)	Petrichloral	
(43)	Phenobarbital	2285*
(44)	Pinazepam	2883-(1-1985)
(45)	Prazepam	
(46)	Quazepam	2881-(11-1986)
(47)	Temazepam	2925-(9-1981)
(48)	Tetrazepam	2886-(1-1985)
(49)	Triazolam	2887-(7-1983)
(50)	Zaleplon	2781-(9-1999)
(51)	Zolpidem	2783-(12-1993)
(52)	Zopiclone	2784-(1-2006)
(53)	Alfaxalone	2731-(2-2014)
(54)	Carisoprodol	8192-(4-1997)
(55)	Tramadol	9752-(8-2007)
(56)	Suvorexant	2223-(8-2014)
(57)	Brexanolone	2400-(4-2021)
(58)	Lemborexant	2245-(4-2021)

(59) Remimazolam	2846-(5-2022)
(60) Daridorexant	2410
(d) <u>Stimulants</u> : Unless specifically excepted or unless listed in another schedule compound, mixture, or preparation which contains any quantity of the follow stimulant effect on the central nervous system, including its salts, isomers (w or geometric), and salts of such isomers whenever the existence of such salts isomers is possible within the specific chemical designation:	ving substances having a hether optical, position,
(1) Cathine ((+)-Norpseudeophedrine	
(2) Diethylpropion	1610*
(3) Fencamfamin	
(4) Fenproporex	
(5) Lorcaserin	
(6) Mazindol	
(7) Mefenorex	1580-(3-1988)
(8) Modafinil	1680-(1-1999)
(9) Pemoline (including organometallic complexes and chelates thereof)	1530*
(10) Phentermine	
(11) Pipradrol	1750-(9-1981)
(12) Serdexmethylphenidate	1729
(13) Sibutramine	1675-(2-1998)
(14) Solriamfetol	1650-(4-2021)
(15) SPA ((-)-1-dimethylamino-1,2,diphenylethane)	1635-(9-1981)
(e) <u>Other substances</u> : Unless specifically excepted or unless listed in another scompound, mixture, or preparation which contains any quantity of the follow including its salts; isomers whether optical, position, or geometric), and salts whenever the existence of such salts, isomers, and salts of isomers is possible.	ving substances, of such isomers,
(1) Pentazocine	9709-(4-1979)
(2) Butorphanol	9720-(4-1997)
(3) Nalbuphine	(4-1997)
(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 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-(3-1988)	
(2) Ephedrine:a -{-(Methylamino)ethyl}benzene-methanol; (10-1995)	
a-{-(methylamino) ethyl} benzyl alcohol; 2-methylamino-1-phenyl-1-propanol; 1-phenyl-1-hydroxy-2-methylaminopropane; 1-phenyl-2-methylaminopropanol; a - hydroxy-b-methylaminopropylbenzene; a product which occurs in the Chinese herb Ma Huang (Ephedra vulgaris, Ephedra sinica Stapf., Ephedra equisetina Bunge, Gnetaceae) in several other Ephedra spp.	
(3) Phenylpropanolamine(7-2005)	

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.

(4) Pseudoephedrine ----- (7-2005)

(e)	material	ants: Unless specifically exempted or excluded or unless listed in another so, compound, mixture, or preparation which contains any quantity of the follow depressant effect on the central nervous system, including its salts, isomers,	wing substances
	(1) Preg	abalin [(S)-3-(aminomethyl)-5-methylhexanoic acid	- 2782-(1-2006)
	(2) Laco	osamide	- 2746-(5-2009)
	(3) Briv	aracetam	- 2710-(4-2017)
		niditan	
		obamate	
	(6) Gana	axolone	2401
(f)	Other s	substances:	
	(1) Non	e.	
<u>SC</u>	HEDUL	E VI ****	
	by a con substance salts, isc Schedule	y from a substance of vegetable origin or independently by means of chemical phination of extraction and chemical synthesis, that contains any quantity of theses, or that contains any of their salts, isomers, and salts of isomers when the owners, and salts of isomers is possible within the specific chemical designation of VI:	the following existence of the in, is included in
	(2) Tetra	ahydrocannabinols, unless the tetrahydrocannabinol is:	**
	(i)	Contained in hemp-derived cannabidiol;	
	(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	(6-2020) ***
	(iii)	Not approved by the United States Food and Drug Administration for marketing as a medication;	(6-2020)
	(3) A sy	nthetic equivalent of:	
	(i)	The substance contained in the Cannabis plant; or	**
	(ii)	The substance contained in the resinous extractives of the genus Cannabis;	**
	class the p	ia divinorum or Salvinorin A, which includes all parts of the plant presently sified botanically as Salvia divinorum, whether growing or not, the seeds of plant, any extract from any part of the plant, and every compound, ufacture, derivative, mixture, or preparation of the plant, its seeds, or its	

	salts, ison	ncluding salts, isomers, and salts of isomers when the existence of the ners, and salts of isomers is possible within the specific chemical on;	**
(5)	classes de specific un Compoun numerical	substances, derivatives, or their isomers in the chemical structural scribed below in subdivisions $(a)(5)(i)-(a)(5)(x)$ of this section and also nclassified substances in subdivision $(a)(5)(xi)$ of this section. ds of the structures described in this subdivision $(a)(5)$ , regardless of designation of atomic positions, are included in this subdivision $(a)(5)$ . etic substances, derivatives, or their isomers included in this subdivision	
	(i) Tetr	ahydrocannabinols:	
	(A) Te	trahydrocannabinols, including without limitation the following:	**
	a)	Delta-1 cis or trans tetrahydrocannabinol [other name(s): Delta-9 cis or trans tetrahydrocannabinol], and its optical isomers;	**
	b)	Delta-6 cis or trans tetrahydrocannabinol [other name(s): Delta-8 cis or trans tetrahydrocannabinol], and its optical isomers;	**
	c)	Delta- 3,4 cis or trans tetrahydrocannabinol [other name(s): Delta-6a,10a cis or trans tetrahydrocannabinol], and its optical isomers;	**
	d)	Delta-10 cis or trans tetrahydrocannabinol, and its optical isomers;	***
	e)	Delta-8 tetrahydrocannabinol acetate ester;	***
	f)	Delta-9 tetrahydrocannabinol acetate ester;	***
	g)	Delta-6a,10a, tetrahydrocannabinol acetate ester;	
	h)	Delta-10 tetrahydrocannabinol acetate ester; and,	***
	i)	A product derived from industrial hemp that was produced as a result of a synthetic chemical process that converted the industrial hemp or a substance contained in industrial hemp into Delta-8, Delta-9, Delta 6a,10a, or Delta-10 tetrahydrocannabinol including their respective acetate esters.	***
	dr Ac	conabinol in sesame oil and encapsulated in a soft gelatin capsule in a aug product approved by the United States Food and Drug dministration is not a tetrahydrocannabinol under this subdivision (5)(i);	**
	na the cy (4- ine rin	hthoylindoles, or any compound structurally derived from 3-(1-phthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at a nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-morpholinyl)ethyl group, whether or not further substituted in the dole ring to any extent and whether or not substituted in the naphthyl ag to any extent, including without limitation the following:	
	(A)JV	/H-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**

(	(B) JWH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	<sup>**</sup>
(	(C) JWH-018, or 1-Propyl-3-(1-naphthoyl)indole;	
(	(D) JWH-019, or 1-Hexyl-3-(1-naphthoyl)indole;	**
	(E) JWH-073, or 1-Butyl-3-(1-naphthoyl)indole;	
(	(F) JWH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
(	(G) JWH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**
(	(H) JWH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl) indole;	**
(	(I) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole;	**
(	(J) JWH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;	**
(	(K) JWH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;	**
(	(L) JWH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole;	**
(	(M) AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole;	**
(	(N) MAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1-naphthalenyl)-methanone;	**
(	(O) EAM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)-methanone; and	**
(	(P) THJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-yl)methanone;	7024-(7-2019)
(iii)	Naphthylmethylindoles, or any compound structurally derived from an H-indol-3-yl-(1-naphthyl) methane by substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation the following:	**
(	(A) JWH-175, or 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane; and	**
(	(B) JWH-184, or 1-Pentyl-1H-3-yl-(4-methyl-1-naphthyl)methane;	**
(iv)	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-	**
(v)	Naphthylmethylindenes, or any compound structurally derived from 1-(1-napthylmethyl)indene with substitution at the 3-position of the indene ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-	

methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indene ring to any extent and whether or not substituted in the naphthyl ring to any extent, including without limitation JWH-176, or E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane;	**
(vi) Phenylacetylindoles, or any compound structurally derived from 3- phenylacetylindole by substitution at the nitrogen atom of the indole ring with alkyl, haloalkyl, alkenyl, cycloalkylmethyl,cycloalkylethyl, 1-(N- methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
(A) JWH-201, or 2-(4-methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone;	**
(B) JWH-203, or 1-Pentyl-3-(2-chlorophenylacetyl)indole;	**
(C) JWH-250, or 1-Pentyl-3-(2-methoxyphenylacetyl)indole;	
(D) JWH-251, or 1-Pentyl-3-(2-methylphenylacetyl)indole; and	**
(E) RCS-8, or 1-(2-cyclohexylethyl)-3-(2- methoxyphenylacetyl)indole;	**
(vii) Cyclohexylphenols, or any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not substituted in the cyclohexyl ring to any extent, including without limitation the following:	**
(A) CP 47,497 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol;	**
(B) Cannabicyclohexanol or CP 47,497 C8 homologue, or 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol; and	**
(C) CP 55,940, or 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-phenol;	**
(viii) Benzoylindoles, or any compound structurally derived from a 3- (benzoyl)indole structure with substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent, including without limitation the following:	**
(A) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole;	**
(B) RCS-4, or 1-Pentyl-3-(4-methoxybenzoyl)indole;	**
(C) WIN-48,098 or Pravadoline, or (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-y]methanone;	**

	(D) AM-2233, or 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole; and	**
	(E) RCS-4 (C4 homologue) or (4-methoxyphenyl)(1-butyl-1H-indol-3-yl)-methanone;	**
(ix)	Adamantoylindoles, or Adamantoylindazoles, including Adamantyl Carboxamide Indoles and Adamantyl Carboxamide Indazoles, or any compound structurally derived from 3-(1-adamantoyl) indole, 3-(1-adamantoyl) indazole, or 3-(2-adamantoyl)indole by substitution at a nitrogen atom of the indole or indazole ring with alkyl, haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl,cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole or indazole ring to any extent and whether or not substituted in the adamantly ring to any extent, including without limitation the following:	**
	(A) AM-1248, or 1-adamantyl-[1-[(1-methylpiperidin-2-yl)methyl]indol-3-yl]methanone;	
	(B) AB-001, or 1-adamantyl-(1-pentylindol-3-yl)methanone;	**
	(C) JWH-018 adamantyl carboxamide, or 1-pentyl-N-tricyclo[3.3.1.13,7]dec-1-yl-1H-indole-3-carboxamide [other name(s): 2NE1, moved in Schedule VI in 2020]	
	(D) AKB-48, or N-(1-adamantyl)-pentyl-1H-indazole-3-carboxamide;	**
	(E) 5F-AKB-48, or N-((3s,5s,7s)-adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	
	(F) STS-135, or N-(1-adamantyl)-1-(5-fluoropentyl)indole-3-carboxamide;	**
(x)	Tetramethylcyclopropylcarbonylindoles or any compound structurally derived from 3-(2,2,3,3-tetramethylcyclopropylcarbonyl) indole by substitution at the nitrogen atom of the indole ring with alkyl,haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl, cycloalkylethyl, (N-methylpiperidin-2-yl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole ring to any extent, including without limitation the following:	**
	(A) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(B) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(C) A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**
	(D) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone;	**

(E) 5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone; and**
(F) A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone;**
(xi) Unclassified Synthetic Cannabinoids, including without limitation the following:**
(A) CP 50556-1 hydrochloride, or [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3- [(2R)-5-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10a- octahydrophenanthridin-1-yl] Acetate;**
(B) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;*
(C) HU-211, or Dexanabinol,(6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;*
(D) Dimethylheptylpyran or DMHP;**
(E) WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone;**
(F) URB597, or [3-(3-carbamoylphenyl)phenyl] N-Cyclohexylcarbamate;*
(G) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one;*
(H) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]-methanone;*
a) URB602, or cyclohexyl N-(3-phenylphenyl)carbamate;**
(I) PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3-carboxylate;**
(J) 5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate;*
(K) BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate;*
(L) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide; **
(M) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3-carboxamide;**
(N) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L-valine methyl ester 7033-(9-2018)
(O) MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)- L-valinate 7044-(9-2018
(P) 5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3-carboxamido)-3,3-dimethylbutanoate; 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) MAB-CHMINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl) 1-(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;	(9-2018)
(Z) 5F-CUMYL-PINACA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7083-(6-2020)
(AA) ADB-FUBINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide	
(BB) 4-Fluoro-MDMB-BUTINACA, or methyl(S)-2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate;	- 7043-(4-2021)
(CC) 5F-AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide	- 7025-(5-2022)
(DD) 4-CN-CUMYL-BUTINACA, or 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide	7089 –(5-2022)
(EE) 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)
(GG) 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate	
(HH) FUB-144, or (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl) methanone;	7014
(II) FUB-AKB48, or N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide;	7047
(JJ)MDMB-4en-PINACA, or Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)butanoate:	-

- (KK) CH-PIATA, or N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide.
- (6) A synthetic substance, derivative, or its isomers with:
  - (i) Similar chemical structure to any substance described in subdivisions (a)(1)(a)(5) of this section; or ------\*\*
  - (ii) Similar pharmacological effects to any substance described in subdivisions (a)(1)-(a)(5) of this section. -----\*\*
- (b) However, except as provided under subsection (c) of this section, the Secretary shall not delete a controlled substance listed in this section from Schedule VI.
- (c) A prescription drug approved by the United States Food and Drug Administration under 21 U.S.C. § 355 is excluded from Schedule VI unless the secretary objects under § 5-64-201.
  - \*-Scheduled before April, 1979.
  - \*\*-Schedule VI is revised to conform to Act 329 of 2013.
  - \*\*\* Schedule VI is revised further to conform to Act 629 of 2023. Each substance added to the Controlled Substances List pursuant to Act 629 of 2023 shall have the following effective dates:
    - (a) For persons who are under twenty-one (21) years of age, the effective date shall be the effective date of Act 629 of 2023; and,
    - (b) For persons who are twenty-one (21) years of age or older, the effective date shall be August 1, 2023.
  - \*\*\*\* Pursuant to ongoing litigation, and a preliminary injunction against enforcing Act 629 of 2023, the changes made to the List of Controlled Substances pursuant to Act 629 of 2023 are not enforceable until a final order issued in the matter, Bio Gen, LLC, et al. v. Sarah Huckabee Sanders, et al., Case No. 4:23-CV-00718-BRW, Central Division, Eastern District of Arkansas, United States District Court.

# QUESTIONNAIRE FOR FILING PROPOSED RULES WITH THE ARKANSAS LEGISLATIVE COUNCIL

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

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

sure to label the markup copy clearly as the markup.

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

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

## FINANCIAL IMPACT STATEMENT

## PLEASE ANSWER ALL QUESTIONS COMPLETELY.

DEP	PARTMENT
	ARD/COMMISSION
PER	SON COMPLETING THIS STATEMENT
TEL	EPHONE NO. EMAIL
emai	omply with Ark. Code Ann. § 25-15-204(e), please complete the Financial Impact Statement and l it 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?

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

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

Yes No

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

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