RECEIVED List **O**f Controlled Substances

FEB 15 2024



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 hereb me and signed the above refe	by certify that Shane David, Pharm.D., well known to me, appeared before erenced document.
Sworn and subscribed	d 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)]	-9815-(2	
(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)]	- 9814-(6	-1982)
(8) Alpha-methylthiofentanyl(N-[1-methyl-2-(2thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	-9832-(2	-1986)
(9)(6) Benzethidine		9606*
(10)(7) Betacetylmethadol		9607*
(11) Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide)]	-9830-(2	-1986)
(12) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2-phenethyl) 3-methyl-4-piperidinyl]-N-phenylpropamamide]		-1986)
(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	813-(10-1985)
(35) 3-methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide)	9833-(2-1986)
(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)-	812 (11-1986)
(42) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2-phenenthyl)-	812-(11-1986)
(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]	812-(11-1986) 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)	812 (11-1986) 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]	-9821 (4-2017)
(57) Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide]	-9822-(4-2017)
(58) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl}-N-phenylpropionamide]	-9836-(4-2017)
(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]	-9840-(4-2017)
(61) Furanyl fentanyl [other name(s): N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylfuran 2-carboxamide]	-9834-(4-2017)
(62) Isobutyryl fentanyl [other name(s): 2-methyl-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide]	-9827-(4-2017)
(63) Ocfentanil [other name(s): N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide]	-9838 (4-2017)
(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]	-9823-(4-2017)
(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)

phenylacrylamide] phenylacrylamide	9811-(6-2020)
(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]	- 9825 (4-2021)
(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)]	
(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
 (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: With or without substitution on the benzimidazole; With or without substitution at the ethylamine; With or without inclusion of the ethylamine in a cyclic structure; With or without substitution on the benzyl ring; or With or without replacement of the benzyl ring with an aromatic ring.
(vi) Benzimidazole-opioid substances shall include but are not limited to:
(A)4'-Hydroxy Nitazene;
(B) 5-Aminoisotonitazene;
(C) Butonitazene;
(D) Clonitazene ;9612 *
(E) Etodesnitazene, [other name(s): Etazene];
(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	``
	(4-2021)
(7) Flualprazolam	(4-2021)
(8) Flubromazepam	(4-2021)
(9) Flubromazolam	(4-2021)
 (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; 	
(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) (<u>+</u>) CIS-4-Methylaminorex [(<u>+</u>)CIS-4,5-dihydro-4- methyl-5-phenyl-2- oxazolamine]	1590-(6-1990)
(iii) Fenethylline 1:	
(iv) N-Benzylpiperazine 74	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>	9- (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	Drugs: 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)	[¹⁸ 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)

⁽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) Depressants : 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) Immediate Precursor : 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.

Administration for marketing only as a suppository.	
(3) Any substance which contains any quantity of a derivative of barbituric acid or any salt thereof	2100*
(4) Chlorhexadol	2510*
(5) Any drug product containing gamma hydroxybutyric acid, including its salts, isomers, and salts of isomers, for which an application is approved under section 505 of the Federal Food, Drug, and Cosmetic Act	
(6) Ketamine. its salts, isomers, and salts of isomers	7285-(7-1999)
Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)-Cyclohexanone.	
(7) Lysergic acid	
(8) Lysergic acid amide	7310*
(9) Methyprylon	2575*
(10) Sulfondiethylmethane	
(11) Sulfonethylmethane	
(12) Sulfonmethane	
(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)
(d) Nalorphine	9400*
(e) Narcotic drugs: Unless specifically excepted or unless listed in another schedule:	7100
(1) Any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:	
(i) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium	9803*
(ii) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9804*

	(ii	i) Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts	9807*
	(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*
	(v	Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts	9810*
		ny material, compound, mixture, or preparation containing any of the following rcotic drugs or their salts, as set forth below:	
	(i)	Buprenorphine (10-02 Transfer)	9064-(6-1985)
	(ii) Reserved	
(f)	composalts,	olic Steroids: Unless specifically excepted or unless listed in another schedule, a bund, mixture, or preparation containing any quantity of the following substance isomers, and salts of isomers whenever the existence of such salts of isomers is pecific chemical designation:	s, including its possible within
	(1) Bo	oldenone;	(9-1991)
	(2) Bo	oldione;	(1-2010)
	(3) Cl	nlorotestosterone (4-chlortestosterone);	(9-1991)
	(4) Cl	ostebol;	(9-1991)
	(5) De	chydrochlormethyltestosterone;	(9-1991)
	(6) D	esoxymethyltestosterone	(1-2010)
	(7) Di	hydrotestosterone (4-dihydrotestosterone);	(9-1991)
	(8) Di	rostanolone;	(9-1991)
	(9) Et	hylestrenol;	(9-1991)
	(10)	Fluoxymesterone;	(9-1991)
	(11)	Formebulone (formebolone);	(9-1991)
	(12)	Mesterolone;	(9-1991)
	(13)	Methandienone, also known as Methandrostenolone;	(9-1991)
	(14)	Methandranone;	(9-1991)
	(15)	Methandriol;	(9-1991)

(16	(i) Methenolone;	(9-1991)
(17	() Methyltestosterone;	(9-1991)
(18	d) Mibolerone;	(9-1991)
(19) Nandrolone;	(9-1991)
(20	19-Nor-4,9(10)-Androstadienedione	(1-2010)
(21) Norethandrolone;	(9-1991)
(22		(
(23	(a) Oxymesterone;	(9-1991)
(24	Oxymetholone;	(9-1991)
(25	s) Stanolone;	(9-1991)
(26	5) Stanozolol;	(9-1991)
(27	() Testolactone;	(9-1991)
(28		
(29) Trenbolone	(9-1991)
(30		(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)
/	empt anabolic steroid products: Compounds, mixtures, or preparations that coroid that have been exempted by the Secretary:	ontain an anabolic
	NDC Number	
(1)	Andro-Estro 90-4	0536-1605
(2)	Androgyn L.A	0456-1005
(3)	Component E-H in Process Pellets	Ivy Labs Inc.
(4)	Component E-H in Process Granulation	Ivy Labs Inc
(5)	Component TE-S in process Granulation	Ivy Labs Inc
(6)	Component TE-S in process Pellets	Ivy Labs Inc
(7)	depANDROGYN	0456-1020
(8)	Depo-Testadiol	0009-0253
(9)	DEPO-T.E	52765-257
(10	depTESTROGEN	51698-257
(11) Duomone	52047-360

(12)	DUO-SPAN II	0684-0102
(13)	DURATESTRIN	43797-016
(14)	Essian	- Pharmaceutics
(15)	Essian H.S	- Pharmaceutics
(16)	Esterified Estrogens & Methyltestosterone, USP (0.625 mg/1.25mg)	Interpharm
(17)	Esterified Estrogens & Methyltestosterone, USP (1.25mg/2.5mg)	Interpharm
(18)	Esterified Estrogens & Methyltestosterone (0.625mg/1.25mg) Tablet	ANDAPharm
(19)	Esterified Estrogens & Methyltestosterone (1.25mg/2.5mg) Tablet	ANDAPharm
(20)	Estratest	0032-1026
(21)	Estratest HS	
(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	
(27)	Premarin with Methyltestosterone	
(28)	Premarin with Methyltestosterone	0046-0878
(29)	Syntest D.S	66576-231
(30)	Stntest H.S	66576-230
(31)	Synovex H in process bulk pellets	
(32)	Synovex H in process granulation	Syntex Animal
(33)	Synovex Plus in process granulation	Fort Dodge
(34)	Synovex Plus in process bulk pellets	Fort Dodge
(35)	TEST-ESTRO Cypionates	0536-9470
(36)	Testoderm with Adhesive 4mg/d	Alza Corp
(37)	Testoderm 4mg/d	17314-4608
(38)	Testoderm 6mg/d	17314-4609
(39)	Testoderm with Adhesive 6mg/d	17314-2836
(40)	Testoderm in process film	Alza Corp
(41)	Testoderm with Adhesive in process film	Alza Corp
(42)	Testosterone Cypionate/Estradiol Cypionate injection	54274-530
(43)	Testosterone Cypionate/Estradiol Cypionate injection	0182-3069
(44)	Testosterone Cyp 50 Estradiol Cyp 2	0814-7737

(45)	Testosterone Cypionate/Estradiol Cypionate injection	0364-6611
(46)	Testosterone Cypionate/Estradiol Cypionate injection	0402-0257
(47)	Testosterone Enanthate/Estradiol Valerate injection	0182-3073
(48)	Testosterone Enanthate/Estradiol Valerate injection	0364-6618
(49)	Testosterone Enanthate/Estradiol Valerate injection	0402-0360
(50)	Testosterone Ophthalmic Solution	Allergan
(51)	Tilapia Sex Reversal Feed (investigational	Ranger, Inc
` /	rinary Anabolic Steroid Implant Products: Anabolic steroid products nistration through implants in cattle or other nonhuman species exemplants	± *
	NDC/DIN	
(1)	Component E-H	021641-002
	Component E-H	
	Component TE-S	
` /	Component T-H	
	Component T-S	
	T-TO	
(7) F	inaplix-H	12799-807-10
` '	inaplix-S	
(9) H	Ieifer-old	
(10)	Heifer-old	
(11)	Heifer-old	
(12)	Implus-H	0009-0434-01
(13)	Implus-H	06-0434-01
	01968327	
(14)	Masculinizing Feed for Fish (Invesitigational)	Rangen,Inc.
(15)	Revalor-G	12799-811
(16)	Revalor-H	12799-810
(17)	Revalor-S	12799-809
(18)	Synovex H	0856-3901
(19)	Synovex H	Syntex
(20)	Synovex Plus	0856-3904
(21)	Tilapia Sex Reversal Feed (investigational)	Zeigier Bros.

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

(i) Hallucinogenic substances:

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

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

SCHEDULE IV

- (a) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.
- (b) <u>Narcotic drugs</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:
 - (1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit------ 9167*
 - (2) Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2-propionoxybutane) ------ 9278-(11-1987)
- (c) <u>Depressants</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) Alprazolam	2882-(6-1982)
(2) Barbital	2145*
(3) Bromazepam	2748-(1-1985)
(4) Camazepam	2749-(1-1985)
(5) Chloral betaine	2460*
(6) Chloral hydrate	2465*
(7) Chlordiazepoxide	2744*
(8) Clobazam	2751-(1-1985)
(9) Clonazepam	2737*
(10) Clorazepate	2768*

(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	
(18)	Ethinamate	
(19)	Ethyl loflazepate	
(20)	Fludiazepam	2759-(1-1985)
(21)	Flunitrazepam	2763-(1-1985)
(22)	Flurazepam	2767*
(23)	Fospropofol	
(24)	Halazepam	2762-(6-1982)
(25)	Haloxazolam	2771-(1-1985)
(26)	Ketazolam	
(27)	Loprazolam	2773-(1-1985)
(28)	Lorazepam	2885*
(29)	Lormetazepam	2774-(1-1985)
(30)	Mebutamate	
(31)	Medazepam	
(32)	Meprobamate	2820*
(33)	Methohexital	2264*
(34)	Methylphenobarbital (mephorbarbital)	2250*
(35)	Midazolam	2884-(1-1985)
(36)	Nimetazepam	2837-(1-1985)
(37)	Nitrazepam	2834-(1-1985)
(38)	Nordiazepam	2838-(1-1985)
(39)	Oxazepam	2835*
(40)	Oxazolam	2839*
(41)	Paraldehyde	2585*
(42)	Petrichloral	2591*
(43)	Phenobarbital	2285*

(44)	Pinazepam	2883-(1-1985)
(45)	Prazepam	2764*
(46)	Quazepam	2881-(11-1986)
(47)	Temazepam	2925-(9-1981)
(48)	Tetrazepam	2886-(1-1985)
(49)	Triazolam	2887-(7-1983)
(50)	Zaleplon	· · · · · · · · · · · · · · · · · · ·
(51)	Zolpidem	2783-(12-1993)
(52)	Zopiclone	
(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
follor of su (1) F (e)(d) comparisments	Iuramine: Any material, compound, mixture, or preparation which wing substances, including its salts, isomers (whether optical, positive isomers, whenever the existence of such salts, isomers, and salts). Senfluramine Stimulants: Unless specifically excepted or unless listed in anotopound, 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	tion, or geometric), and salts sof isomers is possible: 1670* ther schedule, any material, following substances having a ters (whether optical, position,
isom	ers is possible within the specific chemical designation:	
` ′	Cathine ((+)-Norpseudeophedrine	
` /	Diethylpropion	
` /	Gencamfamin	,
	Senproporex	
` /	Lorcaserin	` '
` /	Mazindol	,
(7) N	Mefenorex	1580-(3-1988)

	(8) M	odafinil	1680-(1-1999)
	(9) Pe	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)
(f) (mater substa	Other substances: Unless specifically excepted or unless listed in another schial, compound, mixture, or preparation which contains any quantity of the followances, 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 alts of such
	(1) Pe	entazocineutorphanol	9709-(4-1979)
		albuphine	, ,
	(4) El	luxadoline	9725-(4-2017)
<u>SC</u>	HEDU	ULE V	
(a)		lule V shall consist of the drugs and other substances by whatever official name, name, chemical name, or brand name designated, listed in this section.	common or
(b)	comp	otic Drugs: Unless specifically excepted or unless listed in another schedule, around, mixture or preparation containing any of the following narcotic drugs and rth below.	
	Reser	ved	
(c)	or pre which confe	paration containing any of the following limited quantities of narcotic drugs or so a shall include one or more nonnarcotic active medicinal ingredients in sufficient rupon the compound, mixture, or preparation valuable medicinal qualities other sseed by the narcotic drug alone:	salts thereof, t proportion to
	(1) N	ot more than 200 milligrams of codeine per 100 milliliters or per 100 grams	*
		ot more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 rams.	*
		ot more than 100 milligrams of ethylmorphine per 100 milliliters or per 100 rams.	*

(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)
(4) Pseudoephedrine (7-2005)
Pursuant to Ark. Code Ann. § 5-64-212 as amended in 2005, this Schedule V classification shall NOT apply to any ephedrine, phenylpropanolamine, or pseudoephedrine in liquid, liquid capsule, or liquid gel capsule form. However, sales limits mandated by statute shall apply to all products with ephedrine, phenylpropanolamine, or pseudoephedrine as a listed ingredient regardless of the dosage form.
(e) Depressants: Unless specifically exempted or excluded or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers:
(1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid 2782-(1-2006)
(2) Lacosamide 2746-(5-2009)
(3) Brivaracetam 2710-(4-2017)
(4) Lasmiditan 2790-(4-2021)
(5) Cenobamate 2720-(4-2021)
(6) Ganaxolone 2401
(f) Other substances:

(1) None.

SCHEDULE VI ****

indirectl by a con substance	5-64-214, any material, compound, mixture, or preparation, whether produced day from a substance of vegetable origin or independently by means of chemical subination of extraction and chemical synthesis, that contains any quantity of the ces, or that contains any of their salts, isomers, and salts of isomers when the exi	ynthesis, or following stence of the
salts, iso Schedul	omers, and salts of isomers is possible within the specific chemical designation, is eVI:	is included in
(1) Mar	ijuana	*:
(2) Tetra	ahydrocannabinols, unless the tetrahydrocannabinol is:	*:
(i)	Contained in hemp-derived cannabidiol;	(6-2020
(ii)	Not more than three-tenths of one percent (0.3%) of delta-9 tetrahydrocannabinol in the hemp-derived cannabidiol on a dry weight 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	enthetic 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 man extra salts	ria 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 acts, including salts, isomers, and salts of isomers when the existence of the s, isomers, and salts of isomers is possible within the specific chemical gnation;	*
class spec Com num The	thetic substances, derivatives, or their isomers in the chemical structural sees described below in subdivisions $(a)(5)(i)-(a)(5)(x)$ of this section and also diffic unclassified substances in subdivision $(a)(5)(xi)$ of this section. In appears of the structures described in this subdivision $(a)(5)$, regardless of derivatives designation of atomic positions, are included in this subdivision $(a)(5)$, synthetic substances, derivatives, or their isomers included in this subdivision $(a)(5)$ are:	
(i)	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;	*:

(a) In addition to any substance placed in Schedule VI by the Secretary of the Department of Health

	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.	***
(E	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-chthoyl)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	A)JW	7H-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole;	**
(E	3) JW	7H-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole;	**
(C	C) JW	7H-018, or 1-Propyl-3-(1-naphthoyl)indole;	
(D)JW	7H-019, or 1-Hexyl-3-(1-naphthoyl)indole;	**
(E	E) JW	7H-073, or 1-Butyl-3-(1-naphthoyl)indole;	**
(F) JW	7H-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole;	**
(0	i)JW	7H-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole;	**
(F	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	TH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole;	**
(K	()JW	H-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole;	**
		TH-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;702	24-(7-2019)
(iii)	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 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;	**
y i pinomanono,	

	(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	e; **
(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-carboxyla	
(K) BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxyla	
(L) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxam**	
(M) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole carboxamide;	
(N) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-valine methyl ester	L- 7033-(9-2018)
(O) MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbon L-valinate	nyl)- 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	
(R) MDMB-CHMICA, or methyl 2-(1-(cyclohexylmethyl)-1H- indole-3-carboxamido)-3,3-dimethylbutanoate;	
(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;	
(U) AB-PINACA, or N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H indazole-3-caroboxamide;	
(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 1-(cyclohexylmethyl)-1H-indazole-3-carboxamide;	• /
(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-carboxamide7083-(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 7085-(5-2022
(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 7030
(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; 704
(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.

