

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

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

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

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

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

I, Shane David, Pharm.D., <u>Branch Chief, Health Systems Licensing and Certifications</u> 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-Yates, do hereby certify that Shane David, Pharm.D., well known to me, appeared before me and signed the above referenced document.

Sworn and subscribed to before me this _____ day of March, 2024.

)

Notary Public

My commission expires

ARKANSAS DEPARTMENT OF HEALTH

LIST OF CONTROLLED SUBSTANCES

SECTION I AUTHORITY

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

SECTION II PURPOSE

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

SECTION III GENERAL REQUIREMENTS

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

SECTION IV REPEAL

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

CERTIFICATION

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

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

ARTICLE II

SCHEDULE I

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

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

| (24) | Ketobemidone 96 | 28* |
|------------|---|------|
| (25) | Levomoramide 96 | 29* |
| (26) | Levophenacylmorphan 96 | 31* |
| (27) | Morpheridine 96 | 32* |
| (28) | MPPP [other name(s): (1-methyl-4-phenyl-4-propionoxypiperidine)] 9661-(10-19 | 985) |
| (29) | Noracymethadol 96 | 33* |
| (30) | Norlevorphanol 96 | 34* |
| (31) | Normethadone 96 | 35* |
| (32) | Norpipanone 96 | 36* |
| (33) | PEPAP [other name(s): 1-(2-phenylethyl)-4-phenyl-4 acetyloxypiper-idine] - 9663-(10-19 | 985) |
| (34) | Phenadoxone 96 | 37* |
| (35) | Phenampromide 96 | 38* |
| (36) | Phenomorphan 96 | 47* |
| (37) | Phenoperidine 96 | 41* |
| (38) | Piritramide 96 | 42* |
| (39) | Proheptazine 96 | 43* |
| (40) | Properidine 96 | 44* |
| (41) | Propiram 96 | 49* |
| (42) | Racemoramide 96 | 45* |
| (43) | Tilidine 9750-(9-19 | 981) |
| (44) | Trimeperidine 96 | 46* |
| (45) | Acetyl norfentanyl [other name(s): N-phenyl-N-4-piperidinyl-acetamide](4-20 |)17) |
| (46) di | AH-7921 [other name(s): 3,4-dichloro- <u>N-[(1N-[(1-</u> imethylamino)cyclohexylmethyl]benzamide]9551-(4-20 |)17) |
| | W-18 [other name(s): 1-(4-nitrophenylethyl)piperidylidene-2-(4- hlorophenyl)sulfonamide](4-20 |)17) |
| | W-15 [other name(s): 1-phenylethylpiperidylidene-2-(4- hlorophenyl)Sulfonamide](4-20 |)17) |
| (49) | MT-45 [other name(s): 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine] 9560-(4-20 |)17) |
| (50) N | U-47700 [other name(s): trans-3,4-dichloro-N-(2-(dimethylamino)cyclohexyl)- J-methylbenzamide] 9547-(4-20 |)17) |
| | Fentanyl-related substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Fentanyl-related substance means any substance not therwise listed, and for which no exemption or approval is in effect under section | |

| | 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 [other name(s): N-[1-methyl-2-(2-thienyl)ethyl 4-piperidinyl]-N-phenylpropanamide] | |
| (. | D)Beta-hydroxyfentanyl [other name(s): (N-[1-(2-hydroxy-2-phenethyl)-4- piperidinyl]-N-phenylpropanamide)] | - 9830-(2-1986) |
| (. | E) Beta-hydroxy-3-methylfentanyl [other name(s): N-[1-(2-hydroxy-2- phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropamamide] | - 9831-(2-1986) |
| (| F) 3-Methylfentanyl [other name(s): (N-[3-Methyl-1-(2-phenylethyl)-4- piperidyl]-N-Phenylpropanamide)] | 9813-(10-1985) |
| (| G) 3-methylthiofentanyl [other name(s): N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide] | - 9833-(2-1986) |
| (. | H) Para-fluorofentanyl [other name(s): (N-[4-fluorophenyl)-N-[1-(2- phenenthyl)-4-piperindinyl]propananmide] | 9812-(11-1986) |
| (| I) Thiofentanyl [other name(s): (N-phenyl-N-[1-(2-thienyl)ethyl-4- piperidinyl]-propanamide] | 9835-(2-1986) |
| (. | J) Acetyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N- phenylacetamide] | 9821-(4-2017) |
| (. | K)Butyryl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N- phenylbutyramide] | 9822-(4-2017) |
| (. | L) Beta-hydroxythiofentanyl [other name(s): N-{1-[2-hydroxy-2-(thiophen-2 yl)ethyl]piperidin-4-yl}-N-phenylpropionamide] | |

Page 4

| (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] | |
| (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] | 9842-(5-2022) |

| (FF) 2'-Fluoro ortho-fluorofentanyl [other name(s): N-(1-(2- fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide] 9855-(5-2022) |
|--|
| (GG) 4'-Methyl acetyl fentanyl [other name(s): N-(1-(4-methylphenethyl) piperidin-4-yl)-N-phenylacetamide] 9819-(5-2022) |
| (HH) Ortho-fluorobutyryl fentanyl [other name(s): N-(2-fluorophenyl)-N- (1-phenethylpiperidin-4-yl)butyramide] 9846-(5-2022) |
| (II) Ortho-methyl acetylfentanyl [other name(s): N-(2-methylphenyl)-N-(1- phenethylpiperidin-4-yl)acetamide]9848-(5-2022) |
| (JJ)Ortho-methyl methoxyacetyl fentanyl [other name(s): 2-methoxy-N-(2- methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide] 9820-(5-2022) |
| (KK) Para-methylfentanyl [other name(s): N-(4-methylphenyl)-N-(1- phenethylpiperidin-4-yl)propionamide] 9817-(5-2022) |
| (LL) Phenyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N- phenylbenzamide] 9841-(5-2022) |
| (MM) Thiofuranyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)- N-phenylthiophene-2-carboxamide] 9839-(5-2022) |
| (NN) Fentanyl carbamate [other name(s): ethyl(1-phenethylpiperidin-4- yl)(phenyl)carbamate] 9851-(5-2022) |
| (OO) Ortho-fluoroacryl fentanyl [other name(s): N-(2-fluorophenyl)-N-(1- phenethylpiperidin-4-yl)acrylamide] 9852-(5-2022) |
| (PP) Ortho-fluoroisobutyryl fentanyl [other name(s): N-(2-fluorophenyl)- N-(1-phenethylpiperidin-4-yl)isobutyramide] 9853-(5-2022) |
| (QQ) Para-fluoro furanyl fentanyl [other name(s): N-(4-fluorophenyl)-N- (1-phenethylpiperidin-4-yl)furan-2-carboxamide] 9854-(5-2022) |
| (RR) Meta-fluorofentanyl [other name(s): N-(3-fluorophenyl)-N-(1- phenethylpiperidin-4-yl)propionamide] 9857 |
| (SS) Meta-fluoroisobutyryl fentanyl [other name(s): N-(3-fluorophenyl)- N-(1-phenethylpiperidin-4-yl)isobutyramide9858 |
| (TT) Para-methoxyfuranyl fentanyl [other name(s): N-(4- methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide] 9859 |
| (UU) Para-methylcyclopropyl fentanyl [other name(s): N-(4- methylphenyl)-N-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide] 9865 |
| (VV) 3-furanyl fentanyl [other name(s): N-(1-phenethylpiperidin-4-yl)-N- phenylfuran-3-carboxamide]9860 |
| (WW) 2',5'-dimethoxyfentanyl [other name(s): N-(1-(2,5- dimethoxyphenethyl)piperidin-4-yl)-N-phenylpropionamide] 9861 |
| (XX) Isovaleryl fentanyl [other name(s): 3-methyl-N-(1- phenethylpiperidin-4-yl)-N-phenylbutanamide] 9862 |

Page 6

| | (YY) Ortho-fluorofuranyl fentanyl [other name(s): N-(2-fluorophenyl)-N- (1-phenethylpiperidin-4-yl)furan-2-carboxamide] | 9863 |
|-------------|--|--------------|
| | (ZZ) Alpha'-methyl butyryl fentanyl [other name(s): 2-methyl-N-(1- | |
| | phenethylpiperidin-4-yl)-N-phenylbutanamide] | 9864 |
| (52) | Zipeprol | 9873 |
| (53) | Brorphine | 9098 |
| s a A | Benzimidazole-opioid substances, their isomers, esters, ethers, salts and salts of somers, esters and ethers. Benzimidazole-opioid substances includes any ubstance, not otherwise listed or excepted, and for which no exemption or pproval 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 t its 1-position and benzyl group at its 2-position: | |
| (1 | i) With or without substitution on the benzimidazole; | |
| (1 | ii) With or without substitution at the ethylamine; | |
| (1 | iii) With or without inclusion of the ethylamine in a cyclic structure; | |
| (1 | iv) With or without substitution on the benzyl ring; or | |
| (* | v) With or without replacement of the benzyl ring with an aromatic ring. | |
| (* | vi) Benzimidazole-opioid substances shall include but are not limited to: | |
| | (A)4'-Hydroxy Nitazene; | |
| | (B) 5-Aminoisotonitazene; | |
| | (C) Butonitazene; | |
| | (D)Clonitazene; | 9612 * |
| | (E) Etodesnitazene, [other name(s): Etazene]; | 9765 |
| | (F) Etonitazene ; | 9624* |
| | (G)Flunitazene; | |
| | (H) Isotonitazene; | 9614 |
| | (I) Isotodesnitazene; | |
| | (J) Metodesnitazene; | |
| | (K) Metonitazene; | 9757 |
| | (L) N-Desethyl Etonitazene; | |
| | (M) N-Desethyl Isotonitazene; | |
| | (N)N-Piperidinyl Etonitazene [other name(s): Etonitazepipne]; | |
| | (O) N-Pyrrolidino Etonitazene [other name(s): Etonitazepyne]; | <u> 9758</u> |
| | (P) N-Pyrrolidino Protonitazene; and | |
| | (O) Protonitazene | 9759 |

- (55) <u>2-Methyl AP-237 [other name(s): 1-(2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl)butan-1-one]</u> ------ 9664
- (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* |
| (5) C | odeine-N-Oxide | 9053* |
| (6) C | yprenorphine | 9054* |
| (7) D | esomorphine | 9055* |
| (8) D | ihydromorphine | 9145* |
| (9) D | rotebanol | 9335* |
| (10) | Etorphine (except hydrochloride salt) | 9056* |
| (11) | Heroin | 9200* |
| (12) | Hydromorphinol | 9301* |
| (13) | Methyldesorphine | 9302* |
| (14) | Methyldihydromorphine | 9304* |
| (15) | Morphine methylbromide | 9305* |
| (16) | Morphine methylsulfonate | 9306* |
| (17) | Morphine-N-Oxide | 9307* |
| (18) | Myrophine | 9308* |
| (19) | Nicocodeine | 9309* |
| (20) | Nicomorphine | 9312* |
| (21) | Normorphine | 9313* |
| (22) | Pholcodine | 9314* |
| (23) | Thebacon | 9315* |
| (24) | Mitragynine (11 | -2015) |
| (25) | 7-Hydroxymitragynine(11 | -2015) |

(d) Hallucinogenic substances: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation, which contains any quantity of the following

hallucinogenic substance, or which contains any of its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation (for purposes of this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

| (1) alpha-ethyltryptamine | - 7249-(12-1993) |
|--|------------------|
| Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl)indole; alpha-ET; and AET. | |
| (2) 4-bromo-2,5-dimethoxy-amphetamine | 7391* |
| Some trade or other names: 4-bromo-2,5-dimethoxy-alpha- methylphenethylamine; 4-bromo-2,5-DMA. | |
| (3) 4-bromo-2,5-dimethoxyphenethylamine | 7392-(8-1995) |
| Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1 aminoethan alpha-desmethyl DOB; 2C-B, Nexus. | e; |
| (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 | |

Page 9

| (14) | 3,4,5-trimethoxy amphetamine | 7390* |
|------|--|---------------|
| (15) | 5-methoxy-n,n-dimethyltryptamine 5-MeO-DMT | 7431*(1-2011) |
| (16) | alpha-methyltryptamine | 7432-(7-2005) |
| | Some trade or other names: AMT | |
| (17) | Bufotenine | 7433* |
| | Some trade and other names: 3-(beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine. | |
| (18) | Diethyltryptamine | 7434* |
| | Some trade or other names: N,N-Diethyltryptamine;DET. | |
| (19) | Dimethyltryptamine | 7435* |
| | Some trade or other names: DMT | |
| (20) | 5-methoxy-N,N-diisopropyltryptamine | 7439-(7-2005) |
| | Some trade or other names: 5-MeO-DIPT. | |
| (21) | Ibogaine | 7260* |
| | Some trade and other names: 7-Ethyl-6,6 beta; 7,8,9,10,12,13-octahydro-2- methoxy-6,9-methano-5H-pyrido [1',2': 1,2] azepino [5,4-b] indole; Tabernanthe iboga. | |
| (22) | Lysergic acid diethylamide | 7315* |
| (23) | Mescaline | 7381* |
| (24) | Parahexyl | 7374-(7-1983) |
| | Some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl. | |
| (25) | Peyote | 7415* |
| | Meaning all parts of the plant presently classified botanically as Lophophora williamsii Lemaire, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salts, derivative, mixture or preparation of such plant, its seeds or extracts. (Interprets 21 USC 812 (c), Schedule I (c) (12)). | |
| (26) | N-ethyl-3-piperidyl benzilate | 7482* |
| (27) | N-methyl-3-piperidyl benzilate | 7484* |
| (28) | Psilocybin | 7437* |
| (29) | Psilocyn | 7438* |
| (30) | Ethylamine Analog of phencyclidine | 7455* |

| | Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE. | |
|------------|---|-----------------|
| (31) | Pyrrolidine Analog of phencyclidine | 7458* |
| | Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCPy; PHP | |
| (32) | Thiophene Analog of phencyclidine | 7470* |
| | Some trade or other names: 1-[1-(2-thienyl) cyclohexyl] Piperidine; 2-Thienyl analog of phencyclidine; TPCP; TCP. | |
| (33) | 1-[1-(2-Thienyl)cylcohexyl]pyrrolidine | - 7473-(9-1989) |
| | Some other trade or other names: TCPy. | |
| (34) | N,N-Diallyl-5-Methoxytryptamine; | (6-2012) |
| | Some trade or other names: 5-MeO DALT; 5-Methoxy-DALT | |
| (35) 2: | 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-methoxyphenyl)methyl]ethanamine; 5C-NBOMe | (5-2013) |
| (36) 2: | 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) | 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) 2: | 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 1 | 245-(5-2022) |
| | | Sc | ome trade or other names: Para-methoxymethamphetamine; PMMA | |
| | (54) |) 2- | (ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one | 7286 |
| | | Sc | ome trade or other names: Methoxetamine; MXE | |
| (e) | con dep whe | npour ressa eneve | ants : Unless specifically excepted or unless listed in another schedule, any maind, mixture, or preparation which contains any quantity of the following substant effect on the central nervous system, including its salts, isomers, and salts of er the existence of such salts, isomers, and salts of isomers is possible within the l designation: | nces having a isomers |
| | | hydro and i gamr | ma-hydroxybutyric acid [other name(s): GHB; gamma-hydroxybutyrate; 4- oxybutyrate; 4-hydroxydutanoic acid; sodium oxybate; sodium oxybutyrate], ts known precursors and analogs. Precursors include but are not limited to: ma-butyrolactone2 | · · · · · · · · · · · · · · · · · · · |
| | (2) | Mec | loqualone | 2572* |
| | (3) | Meth | naqualone | 2565* |
| | | esters other under that s pheny on th any c | codiazepine substances, their isomers, esters, ethers, salts and salts of isomers, s and ethers. Benzodiazepine substances includes any substance, not rwise listed or excepted, and for which no exemption or approval is in effect r section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355], structurally has a fused 1,4-diazepine and benzene ring structure with a yl connected to the diazepine ring, with any substitution(s) or replacement(s) the 1,4-diazepine or benzene ring, any substitution(s) on the phenyl ring, or combination thereof. Benzodiazepine substances shall include but are not ed 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]. | |
| | | isomonot o effec | nodiazepine substances, their isomers, esters, ethers, salts and salts of ers, esters and ethers. Thienodiazepine substances includes any substance, otherwise listed or excepted, and for which no exemption or approval is in t under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. , 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) | 2-2 | .01 | 4 |) |
|-----|----------|------|-----|-----|---|---|
|-----|----------|------|-----|-----|---|---|

(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) |
|--|-----------------|
| (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 [other name(s): BZP, 1-Benzylpiperazine] | - 7493-(1-2005) |
| Some trade or other name: BZP, 1-Benzylpiperazine | |
| (v) N-ethylamphetamine | - 1475-(6-1982) |
| (vi) N-[1-(1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers | 9818-(2-1986) |
| (vii) N-[1-(2-thienyl)methyl-4-piperidyl)-N-phenylpropan-mide (thenylfentanyl its optical isomers, salts and salts of isomers | |
| (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 <u>[other name(s):(some</u> -2-Methylamine-Proprophenone, alph (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 | I- |
| (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) |
| (xi) 4,4'-Dimethylaminorex –[other name(s): some 4,4'-DMAR, 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine, or 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine] | 1595-(5-2022) |
| (xii) Amineptine | 1219 |
| (xiii) Mesocarb | 1227 |
| (xiv) Methyl-N-ethylcathinone (MEC) | (6-2014) |
| | |

| | (xv) | Methiopropamine. [other name(s): N-methyl-1-(thiophen-2-yl)propan-2- amine] | 1478 |
|-----|--|--|-----------------|
| (2) | indire 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 ains 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 fife chemical designation, with the following chemical structure is included i dule I: | ; |
| | (i) | 4-Methylmethcathinone (Mephedrone) | - 1248-(3-2011) |
| | (ii) | Methylenedioxypyrovalerone (MDPV) | (3-2011) |
| | (iii) | 3,4-Methylenedioxy-N-methylcathinone (Methylone) | - 7540-(3-2011) |
| | (iv) | 4-Methoxymethcathinone | (3-2011) |
| | (v) | 3-Fluoromethcathinone | (3-2011) |
| | (vi) | 4-Fluoromethcathinone | (3-2011) |
| | (vii) | 1-(1,3-benzodioxol-5-yl)-2-methylamino)butan-1-one-(Butylone) | 7541-(11-2014) |
| | (viii) | Alpha-Pyrrolidinopentiophenone (Alpha-PVP) | 7545-(11-2015) |
| | (ix) | 4-methyl-N-ethylcathinone (4-MEC) | 1249-(9-2018) |
| | (x) | 4-methyl-alpha-pyrrolidinopropiophenone (4-MePPP) | 7498-(9-2018) |
| | (xi) | 2-(methylamino)-1-phenylpentan-1-one (Pentedrone) | 1246-(9-2018) |
| | (xii) | 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Pentylone; bk- MBDP) | 7542-(9-2018) |
| | (xiii) | 4-fluoro-N-methylcathinone (4-FMC, Flephedrone) | 1238-(9-2018) |
| | (xiv) | 3-fluoro-N-methylcathinone (3-FMC) | 1233-(9-2018) |
| | (xv) | 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Naphyrone) | 1258-(9-2018) |
| | (xvi) | Alpha-pyrrolidinobutiophenone ([Alpha]-PBP) | 7546-(9-2018) |
| | <u>(xvii</u> |)3-methylmethcathinone (3–MMC) | 1259 |
| | (xvii | (xviii) | A |
| | | compound, unless listed in another schedule or a legend drug, that is structurally derived from 2-Amino-phenyl-1-propanone by modification of by substitution: | or |
| | | 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 | D) At the 2 position with an alloyd substituents on | |

(B) At the 3-position with an alkyl substituent; or

(C) At the nitrogen atom with alkyl or dialkyl groups, or by inclusion of the nitrogen atom in a cyclic structure.

| (xviii)(xix) |] |
|---|---------------|
| (xviii)(xix) | /543-(7-2019) |
| (xix)(xx) | 1 |
| -(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (Ethylone)7 | 7547-(4-2021) |
| (xx)(xxi) | 1 |
| -(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (Eutylone) | /549-(4-2021) |
| (xxi)(xxii) | 2 |
| -(ethylamino)-1-phenylhexan-1-one [other name(s): (N-Ethylhexedrone ; Alpha-Ethylaminohexanophenone)] | |
| (xxii)(xxiii) | |
| 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one [other name(s): (Alpha- Pyrrolidinohexanophenone ; Alpha-PHP)] | |
| (xxiii)(xxiv) | |
| (xxiii)(xxiv) | 7245 |
| (xxiv)(xxv) | |
| (xxiv)(xxv) | 7446 |
| (xxv)(xxvi) | |
| -phenyl-2-(pyrrolidin-1-yl)heptan-1-one [other name(s): (Alpha- Pyrrolidinoheptaphenone ; PV8)] | |
| (xxvi)(xxvii) | |
| 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): (4'- Chloro-alpha-pyrrolidinovalerophenone ; 4-chloro-alpha-PVP)] | 7443 |
| (xxvii)(xxviii) | |
| 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one [other name(s): alpha- PiHP, Alpha-Pyrrolidinoisohexanophenone] | |

SCHEDULE II

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

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

| | (i) | Raw opium | 9600* |
|-----|-------------------------|--|---------------|
| | (ii) | Opium extracts | 9610* |
| | (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 | 9150* |
| | (xiii) | Metopon | 9260* |
| | (xiv) | Morphine | 9300* |
| | (xv) | Oripavine | 9330*(9-2007) |
| | (xvi) | Oxycodone | 9143* |
| | (xvii |)Oxymorphone | |
| | (xvii | i) Thebaine | 9333* |
| | (xix) | Tapentadol | 9780-(5-2009) |
| | (xx) | Noroxymorphone | 9668-(4-2021) |
| (2) | equiv of the | 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 | m poppy and poppy straw.* | |
| (4) | leave deriv deriv | leaves (9040) and any salt, compound, derivative, or preparation of coca es, (including cocaine (9041) and ecgonine (9180) and their salts, isomers, atives and salts of isomers and derivatives), and any salt, compound, ative, or preparation thereof which is chemically equivalent or identical with of these substances, except that the substances shall not include: | * |
| | (i) | Decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine ; | |
| | (::) | | |

(ii) [¹²³I]ioflupane; or

(iii) $[^{18}F]FP-CIT.$

| (5) Concentrate or poppy straw (the crude extract of poppy straw in either liquid, |
|--|
| solid or powder form which contains the phenanthrene alkaloids of the opium |
| poppy),9670.* |

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

| (1) A | lfentanil | 9737-(2-1987) |
|------------|---|----------------|
| (2) A | lphaprodine | 9010* |
| (3) A | nileridine | 9020* |
| (4) B | ezitramide | 9800* |
| (5) B | ulk Dextropropoxyphene (non-dosage forms) | 9273-(9-1981) |
| (6) C | arfentanil | 9743-(9-1988) |
| (7) D | ihydrocodeine | |
| (8) D | iphenoxylate | |
| (9) Fe | entanyl | 9801* |
| (10) | Isomethadone | |
| (11) | Levo-alphacetylmethadol (LAAM) | 9648-(12-1993) |
| (12) | Levomethorphan | |
| (13) | Levorphanol | |
| (14) | Metazocine | |
| (15) | Methadone | |
| (16) | Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane | 9254* |
| (17) ca | Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane- rboxylic acid | 9802* |
| (18) | Pethidine (Meperidine) | |
| (19) | Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine | |
| (20) | Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate | |
| (21) | Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid | |
| (22) | Phenazocine | |
| (23) | Piminodine | |
| (24) | Racemethorphan | |
| (25) | Racemorphan | |

| (26) | Remifentanil | 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 | 1105* |
| (3) Lisdexamefetamine | 1205*(7-2007) |
| (4) Phenmetrazine and its salts | |
| (5) Methylphenidate | 1724* |

(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) Amobarbital | | 2125* |
|-------------------|---------|--------|
| (2) Glutethimide | 2550-(2 | -1991) |
| (3) Pentobarbital | | 2270* |
| (4) Phencyclidine | e | 7471* |
| (5) Secobarbital | | 2315* |

(f) Hallucinogenic Substances:

- (1) Nabilone ------ 7379-(11-1987)
 [Other name(s) for nabilone: (±)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one].
 (2) Dronabinol in an oral solution in a drug product approved for marketing by the
 Use the set of the se
 - U.S. Food and Drug Administration; [(-)-delta-9-transtetrahydrocannabinol(delta-9-THC)].-----7365-(7-2019)
- (g) **Immediate Precursor**: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

- (1) Immediate precursor to Amphetamine and Methamphetamine:
- (i) Phenylacetone ------ 8501-(3-1980) 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:

| (i) | Amobarbital | 2126* |
|------|--------------|-------|
| (ii) | Secobarbital | 2316* |

| (i | ii) Pentobarbital | 2271* |
|--------------------------|---|-----------------|
| (i | v) Embutramide | 2020*(9-2006) |
| | or any salt thereof and one or more other active medicinal ingredients which are not listed in any schedule. | |
| (2) A | ny suppository dosage form containing: | |
| (i |) Amobarbital | 2126* |
| (i | i) Secobarbital | 2316* |
| (i | ii) Pentobarbital | 2271* |
| | or any salt of any of these drugs and approved by the Food and Drug Administration for marketing only as a suppository. | |
| | ny substance which contains any quantity of a derivative of barbituric acid or | |
| | ny salt thereof | |
| (4) C | hlorhexadol | 2510* |
| | ny drug product containing gamma hydroxybutyric acid, including its salts, | |
| | omers, and salts of isomers, for which an application is approved under section 05 of the Federal Food, Drug, and Cosmetic Act | - 2012-(2-2001) |
| | etamine. its salts, isomers, and salts of isomers | × / |
| | Some other names for Ketamine: (+-)-2-(2-Chlorophenyl)-2-(Methylamino)- Cyclohexanone. | () |
| (7) L | ysergic acid | 7300* |
| (8) L | ysergic acid amide | 7310* |
| | fethyprylon | |
| (10) | Sulfondiethylmethane | |
| (11) | Sulfonethylmethane | 2605* |
| (12) | Sulfonmethane | 2610* |
| (13) | Tiletamine and zolazepam or any salt thereof | - 7295-(3-1988) |
| | Some trade or other name for a tiletamine- zolazepam combination product: Telazol. | |
| | Some trade or other names for tiletamine:2_(ethylamino)-2-(2-thienyl)-cyclohexanone. | |
| | Some trade or other names for zolazepam:4_(2-fluorophenyl)-6,8-dihydro-1,3,8,-trimethyl_pyrazolo-[3,4-e]-[1,4,]-diazepin-7(1-H)-one-; flupyrazapon. | |
| (14) | Perampanel | 2261-(11-2013) |
| <u>(15)</u> <u>ar</u> | Xylazine and any material, compound, mixture, or preparation which contains ny quantity of xylazine, including its salts, isomers, and salts of isomers | |

whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation, except in the following uses:

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

(A) By a licensed pharmacist or licensed veterinarian ; or

- (B) Pursuant to a valid prescription from a licensed veterinarian;
- (iv) Possessing, manufacturing, distributing, or using xylazine as an active pharmaceutical ingredient for manufacturing an animal drug either:
 - (A) Approved under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b); or
 - (B) Issued an investigation use exemption under section 512 of the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 360b(j));
- (v) Manufacturing, distributing, or using a xylazine bulk chemical for pharmaceutical compounding by a licensed pharmacist or veterinarian; or

(vi) Another use approved or permissible under the Federal Food, Drug and Cosmetic Act (21 U.S.C. § 301, et seq.) or under 21 CFR Part 530, Subpart B.

| (d) Nalorphine | 9400* |
|----------------|-------|
|----------------|-------|

(e) **Narcotic drugs:** Unless specifically excepted or unless listed in another schedule:

- Any material, compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below:
 - (i) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium ------ 9803*
 (ii) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic
 - ingredients in recognized therapeutic amounts ------ 9804*
 (iii) Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active nonnarcotic ingredients in recognized therapeutic amounts ------ 9807*

| | (i | 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* |
|-----|----------------|--|-----------------------------------|
| | (1 | 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* |
| | (1 | 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* |
| | | Any material, compound, mixture, or preparation containing any of the following arcotic drugs or their salts, as set forth below: | |
| | (i | i) Buprenorphine (10-02 Transfer) | 9064-(6-1985) |
| | (i | ii) Reserved | |
| (f) | comp salts, | polic Steroids: Unless specifically excepted or unless listed in another schedule, a pound, mixture, or preparation containing any quantity of the following substance isomers, and salts of isomers whenever the existence of such salts of isomers is p pecific chemical designation: | es, including its possible within |
| | (1) B | Boldenone; | (9-1991) |
| | (2) B | Boldione; | (1-2010) |
| | (3) C | Chlorotestosterone (4-chlortestosterone); | (9-1991) |
| | (4) C | Clostebol; | (9-1991) |
| | (5) D | Dehydrochlormethyltestosterone; | (9-1991) |
| | (6) D | Desoxymethyltestosterone | (1-2010) |
| | (7) D | Dihydrotestosterone (4-dihydrotestosterone); | (9-1991) |
| | (8) D | Drostanolone; | (9-1991) |
| | (9) E | Sthylestrenol; | (9-1991) |
| | (10) | Fluoxymesterone; | (9-1991) |
| | (11) | Formebulone (formebolone); | |
| | (12) | Mesterolone; | (9-1991) |
| | (13) | Methandienone, also known as Methandrostenolone; | (9-1991) |
| | (14) | Methandranone; | (9-1991) |
| | (15) | Methandriol; | (9-1991) |
| | (16) | Methenolone; | (9-1991) |
| | (17) | Methyltestosterone; | (9-1991) |

| (18) | Mibolerone; | (9-1991) |
|---|---|--|
| (19) | Nandrolone; | (9-1991) |
| (20) | 19-Nor-4,9(10)-Androstadienedione | (1-2010) |
| (21) | Norethandrolone; | (9-1991) |
| (22) | Oxandrolone; | (9-1991) |
| (23) | Oxymesterone; | (9-1991) |
| (24) | Oxymetholone; | (9-1991) |
| (25) | Stanolone; | (9-1991) |
| (26) | Stanozolol; | (9-1991) |
| (27) | Testolactone; | (9-1991) |
| (28) | Testosterone; | (9-1991) |
| (29) | Trenbolone | (9-1991) |
| (30) | Prostanozol | (8-2012) |
| (31) | Methasterone | (8-2012); |
| | and | |
| (22) | Annualt action an increase of a dama on substance described on list in this | |
| (32) p | Any salt, ester, or isomer of a drug or substance described or list in this aragraph, if that salt, ester, or isomer promotes muscle growth | (9-1991) |
| p (g) Exen | | |
| p (g) Exen | aragraph, if that salt, ester, or isomer promotes muscle growth | |
| p (g) Exen stero | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic |
| p (g) Exen stero (1) A | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic |
| p (g) Exen stero (1) A (2) A | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic |
| p (g) Exen stero (1) A (2) A (3) C | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. |
| p (g) Exen stero (1) A (2) A (3) C (4) C | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc |
| p (g) Exenstero (1) A (2) A (3) C (4) C (5) C | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d (8) E | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc 0456-1020 0009-0253 |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d (8) E | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc 0456-1020 0456-1020 0456-1020 52765-257 |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d (8) D (9) D | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc Ivy Labs Inc 0456-1020 0456-1020 52765-257 51698-257 |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d (8) E (9) E (10) | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc 0456-1020 0456-1020 0456-1020 52765-257 52047-360 |
| p (g) Exensistero (1) A (2) A (3) C (4) C (5) C (6) C (7) d (8) E (9) E (10) (11) | aragraph, if that salt, ester, or isomer promotes muscle growth | contain an anabolic 0536-1605 0456-1005 Ivy Labs Inc. Ivy Labs Inc Ivy Labs Inc 0456-1020 0456-1020 0009-0253 52765-257 51698-257 52047-360 0684-0102 |

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

| (47) | Testosterone Enanthate/Estradiol Valerate injection | - 0182-3073 |
|------|---|-------------|
| (48) | Testosterone Enanthate/Estradiol Valerate injection | - 0364-6618 |
| (49) | Testosterone Enanthate/Estradiol Valerate injection | - 0402-0360 |
| (50) | Testosterone Ophthalmic Solution | Allergan |
| (51) | Tilapia Sex Reversal Feed (investigational) | Ranger, Inc |

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

| N | DC | D | IN |
|---|----|---|----|
| | | | |

| (1) Component E-H | |
|---|---------------|
| (2) Component E-H | 01968327 |
| (3) Component TE-S | 021641-004 |
| (4) Component T-H | 0211641-006 |
| (5) Component T-S | 0211641-005 |
| (6) F-TO | 00093351 |
| (7) Finaplix-H | 12799-807-10 |
| (8) Finaplix-S | 12799-807-07 |
| (9) Heifer-old | Boehringer |
| (10) Heifer-old | Ingelheim |
| (11) Heifer-old | Ivy Lab. |
| (12) Implus-H | 0009-0434-01 |
| (13) Implus-H | 06-0434-01 |
| 01968327 | |
| (14) Masculinizing Feed for Fish (Invesitigational) | Rangen,Inc. |
| (15) Revalor-G | 12799-811 |
| (16) Revalor-H | 12799-810 |
| (17) Revalor-S | 12799-809 |
| (18) Synovex H | 0856-3901 |
| (19) Synovex H | Syntex |
| (20) Synovex Plus | 0856-3904 |
| (21) Tilapia Sex Reversal Feed (investigational) | Zeigier Bros. |
| | |

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

(i) Hallucinogenic substances:

 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)

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[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]
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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:
 - Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit------ 9167*
 Dextro propoxyphene (alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-methyl-2propionoxybutane) ------ 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 | |
|------|--------------------------------------|----------------|
| (15) | Dichloralphenazone | 2467-(10-2002) |
| (16) | Estazolam | 2756-(1-1985) |
| (17) | Ethchlorvynol | |
| (18) | Ethinamate | |
| (19) | Ethyl loflazepate | |
| (20) | Fludiazepam | |
| (21) | Flunitrazepam | |
| (22) | Flurazepam | 2767* |
| (23) | Fospropofol | 2138-(11-2009) |
| (24) | Halazepam | |
| (25) | Haloxazolam | 2771-(1-1985) |
| (26) | Ketazolam | |
| (27) | Loprazolam | |
| (28) | Lorazepam | 2885* |
| (29) | Lormetazepam | |
| (30) | Mebutamate | 2800* |
| (31) | Medazepam | 2836-(1-1985) |
| (32) | Meprobamate | 2820* |
| (33) | Methohexital | 2264* |
| (34) | Methylphenobarbital (mephorbarbital) | 2250* |
| (35) | Midazolam | 2884-(1-1985) |
| (36) | Nimetazepam | 2837-(1-1985) |
| (37) | Nitrazepam | 2834-(1-1985) |
| (38) | Nordiazepam | 2838-(1-1985) |
| (39) | Oxazepam | 2835* |
| (40) | Oxazolam | 2839* |
| (41) | Paraldehyde | 2585* |
| (42) | Petrichloral | 2591* |
| (43) | Phenobarbital | 2285* |
| (44) | Pinazepam | 2883-(1-1985) |
| (45) | Prazepam | 2764* |
| (46) | Quazepam | 2881-(11-1986) |

| (47) | Temazepam | 2925-(9-1981) |
|------|--------------|----------------|
| (48) | Tetrazepam | 2886-(1-1985) |
| (49) | Triazolam | 2887-(7-1983) |
| (50) | Zaleplon | 2781-(9-1999) |
| (51) | Zolpidem | 2783-(12-1993) |
| (52) | Zopiclone | 2784-(1-2006) |
| (53) | Alfaxalone | 2731-(2-2014) |
| (54) | Carisoprodol | 8192-(4-1997) |
| (55) | Tramadol | 9752-(8-2007) |
| (56) | Suvorexant | 2223-(8-2014) |
| (57) | Brexanolone | 2400-(4-2021) |
| (58) | Lemborexant | 2245-(4-2021) |
| (59) | Remimazolam | 2846-(5-2022) |
| (60) | Daridorexant | 2410 |
| (61) | Zuranolone | 2420 |

(d) <u>Stimulants</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers (whether optical, position, or geometric), and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

| (1) Ca | athine ((+)- <u>Norpseudoephedrine)</u> Norpseudeophedrine | 1230-(3-1988) |
|--------|---|---------------|
| (2) D | iethylpropion | 1610* |
| (3) Fe | encamfamin | 1760-(3-1988) |
| (4) Fe | enproporex | 1575-(3-1988) |
| (5) Lo | prcaserin | 1625-(6-2013) |
| (6) M | azindol | 1605-(6-1982) |
| (7) M | efenorex | 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 | 1650-(4-2021) |
|------|--|---------------|
| (15) | SPA ((-)-1-dimethylamino-1,2,diphenylethane) | 1635-(9-1981) |

(e) <u>Other substances</u>: Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts; isomers whether optical, position, or geometric), and salts of such isomers, whenever the existence of such salts, isomers, and salts of isomers is possible:

| (1) Pentazocine | |
|-----------------|--|
| (2) Butorphanol | |
| (3) Nalbuphine | |
| (4) Eluxadoline | |

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 |
| (4) Pseudoephedrine (7-2005 |

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

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

| (1) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid | - 2782-(1-2006) |
|---|-----------------|
| (2) Lacosamide | - 2746-(5-2009) |
| (3) Brivaracetam | - 2710-(4-2017) |
| (4) Lasmiditan | - 2790-(4-2021) |
| (5) Cenobamate | - 2720-(4-2021) |
| (6) Ganaxolone | 2401 |

(f) Other substances:

(1) None.

SCHEDULE VI ****

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

| (1) Marijuana | ** |
|---------------|----|
|---------------|----|

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

| | i) | A product derived from industrial hemp that was produced as a result of a synthetic chemical process that converted the industrial hemp or a substance contained in industrial hemp into Delta-8, Delta-9, Delta 6a,10a, or Delta-10 tetrahydrocannabinol including their respective acetate esters. | *** |
|-------|---------------------------------|---|-----------|
| | drı Ad | ronabinol in sesame oil and encapsulated in a soft gelatin capsule in a ug product approved by the United States Food and Drug dministration is not a tetrahydrocannabinol under this subdivision 0(5)(i); | ** |
| (ii) | naj the cyc (4- inc | hthoylindoles, or any compound structurally derived from 3-(1- phthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at e 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 ng to any extent, including without limitation the following: | ** |
| | (A)JW | VH-007, or 1-pentyl-2-methyl-3-(1-naphthoyl)indole; | ** |
| | (B) JW | VH-015, or 1-Propyl-2-methyl-3-(1-naphthoyl)indole; | ** |
| | | VH-018, or 1-Propyl-3-(1-naphthoyl)indole; | |
| | (D) JW | VH-019, or 1-Hexyl-3-(1-naphthoyl)indole; | ** |
| | (E) JW | VH-073, or 1-Butyl-3-(1-naphthoyl)indole; | ** |
| | (F) JW | VH-081, or 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole; | ** |
| | (G)JW | VH-098, or 1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole; | ** |
| | (H)JW | VH-122, or 1-Pentyl-3-(4-methyl-1-naphthoyl)indole; | ** |
| | (I) JW | VH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole; | ** |
| | (J) JW | VH-200, or 1-[2-(4-morpholiny)ethyl]-3-(1-naphthoyl) indole; | ** |
| | (K)JW | VH-210, or 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole; | ** |
| | (L) JW | VH-398, or 1-Pentyl-3-(4-chloro-1-naphthoyl)indole; | ** |
| | (M) | AM-2201, or 1-(5-fluoropentyl)-3-(1-naphthoyl)indole; | ** |
| | | AM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-methyl-1- phthalenyl)-methanone; | ** |
| | | AM2201, or (1-(5-fluoropentyl)-1H-indol-3-yl)(4-ethyl-1-naphthalenyl)- ethanone; and | ** |
| | · / | HJ-2201, or [1-(5-fluoropentyl)-1H-indazol-3-yl](naphthalen-1-)methanone;7024 | -(7-2019) |
| (iii) | - | hthylmethylindoles, or any compound structurally derived from an H- dol-3-yl-(1-naphthyl) methane by substitution at the nitrogen atom of the | |

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 | ** |
| (1 | 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; | ** |
| (1 | B) JWH-203, or 1-Pentyl-3-(2-chlorophenylacetyl)indole; | ** |
| (| C) JWH-250, or 1-Pentyl-3-(2-methoxyphenylacetyl)indole; | ** |
| (1 | 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; | ** |
| | phonon, | |

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

| (x) | Tetramethylcyclopropylcarbonylindoles or any compound structurally derived from 3-(2,2,3,3-tetramethylcyclopropylcarbonyl) indole by substitution at the nitrogen atom of the indole ring with alkyl,haloalkyl, alkenyl, cyanoalkyl, hydroxyalkyl, cycloalkylmethyl, cycloalkylethyl, (N- methylpiperidin-2-yl)methyl or 2-(4-morpholinyl)ethyl, whether or not further substituted in the indole ring to any extent, including without limitation the following: | ** |
|------|---|----|
| | (A) UR-144, or (1-pentylindol-3-yl)-(2,2,3,3- tetramethylcyclopropyl)methanone; | ** |
| | (B) XLR-11, or [1-(5-fluoropentyl)-1H-indol-3-yl]-(2,2,3,3- tetramethylcyclopropyl)methanone; | ** |
| | (C) A-796,260, or [1-(2-morpholin-4-yl-ethyl)-1H-indol-3-yl](2,2,3,3- tetramethylcyclopropyl)methanone; | ** |
| | (D) 5-Chloro-UR-144, or ([-(5-chloropentyl)-1H-indol-3-yl](2,2,3,3- tetramethylcyclopropyl)methanone; | ** |
| | (E) 5-Bromo-UR-144, or [1-(5-bromopentyl)-1H-indol-3-yl](2,2,3,3- tetramethylcyclopropyl)methanone; and | ** |
| | (F) A-834,735, or 1-(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3- tetramethylcyclopropyl)methanone; | ** |
| (xi) |) Unclassified Synthetic Cannabinoids, including without limitation the following: | ** |
| | (A) CP 50556-1 hydrochloride, or [(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3- [(2R)-5-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10a- octahydrophenanthridin-1-yl] Acetate; | ** |
| | (B) HU-210, or (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2- methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol; | ** |
| | (C) HU-211, or Dexanabinol,(6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3- (2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol; | ** |
| | (D) Dimethylheptylpyran or DMHP; | ** |
| | (E) WIN55,212-2, or 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl-1-naphthalenylmethanone; | ** |
| | (F) URB597, or [3-(3-carbamoylphenyl)phenyl] N-Cyclohexylcarbamate; | ** |
| | (G) URB754, or 6-methyl-2-[(4-methylphenyl)amino]-1-benzoxazin-4-one; | ** |
| | (H) CB-13, or 1-naphthalenyl[4-(pentyloxy)-1 naphthalenyl]-methanone; | ** |
| | a) URB602, or cyclohexyl N-(3-phenylphenyl)carbamate; | ** |
| | (I) PB-22, or quinolin-8-yl 1-(5-pentyl)-1H-indole-3-carboxylate; | ** |
| | (J) 5F-PB-22, or quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate; | ** |

| (K) BB-22, or quinolin-8-yl 1-(cyclohexylmethyl)-1H-indole-3-carboxylate; ** |
|--|
| (L) NNEI (MN-24), or N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide; ** |
| (M) 5F-NNEI, or 1-(5-fluoropentyl)-N-(naphthalene-1-yl)-1H-indole-3- carboxamide; ** |
| (N) 5-Fluoro-AMB, or n-[[1-(5-fluoropentyl)-1H-indazol-3-yl]carbonyl]-L- valine methyl ester 7033-(9-2018) |
| (O) MMB-CHMICA,or methyl-(1-cyclohexylmethyl)-1H-indole-3-carbonyl)- L-valinate 7044-(9-2018) |
| (P) 5-Fluoro-ADB, or methyl 2-(1-(5-fluoropentyl)-1H- indazole-3- carboxamido)-3,3-dimethylbutanoate; 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-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) |

| | (E | E) | 5F-CUMYL-P7AICA, or 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)- 1H-pyrrolo[2,3-b]pyridine-3-carboxamide | 7085-(5-2022) |
|-----|-----------|-----------|--|---------------|
| | (F | F) | NM2201, or Naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3- carboxylate | 7221-(5-2022) |
| | (G | ίG) | 5F-EDMB-PINACA, or Ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3- carboxamido)-3,3-dimethylbutanoate | 7036 |
| | (Н | IH) | 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; | |
| | (JJ | J) | MDMB-4en-PINACA, or Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)- 1H-indazole-3-carboxamido)butanoate; | |
| | (K | K) | CH-PIATA, or N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide; | |
| | <u>(L</u> | <u>L)</u> | ADB-BUTINACA, or N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1- butyl-1H-indazole-3-carboxamide | 7027 |
| (6) | A syn | the | tic substance, derivative, or its isomers with: | |
| | (i) S | | nilar chemical structure to any substance described in subdivisions (a)(1)- a)(5) of this section; or | |
| | (ii) S | | nilar 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.