

**Title 19 - DEPARTMENT OF HEALTH AND  
SENIOR SERVICES  
Division 30—Division of Regulation and Licensure  
Chapter 1—Controlled Substances**

**PROPOSED AMENDMENT**

**19 CSR 30-1.002 Schedules of Controlled Substances.** The department is amending section (1).

*PURPOSE: This amendment updates the Schedules of Controlled Substances to be consistent with 21 CFR Part 1308.*

*PUBLISHER'S NOTE: The secretary of state has determined that the publication of the entire text of the material which is incorporated by reference as a portion of this rule would be unduly cumbersome or expensive. This material as incorporated by reference in this rule shall be maintained by the agency at its headquarters and shall be made available to the public for inspection and copying at no more than the actual cost of reproduction. This note applies only to the reference material. The entire text of the rule is printed here.*

(1) Schedules of Controlled Substances.

(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 Drug Enforcement Administration (DEA) Controlled Substances Code Number set forth opposite it.

1. Opiates. 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, and salts is possible within the specific chemical designation:

- |  |      |
|--|------|
| A. Acetyl-alpha-methylfentanyl<br>(N-(1-(1-methyl-2-phenethyl)-<br>4-piperidinyl)-N-<br>phenylacetamide)   | 9815 |
| B. Acetylmethadol  | 9601 |
| C. Acetyl fentanyl (N-(1-<br>phenethylpiperidin-4-yl)-<br>N-phenylacetamide)   | 9821 |
| D. N-(1-phenethylpiperidin-<br>4-yl)-N-phenylacrylamide,<br>its isomers, esters, ethers,<br>salts, and salts of isomers,<br>esters, and ethers (other<br>names: acryl fentanyl,<br>acryloylfentanyl) | 9811 |
| E. AH-7921(3,4-dichloro-<br>N-[(1-dimethylamino)<br>cyclohexylmethyl]<br>benzamide)  | 9551 |
| F. Allylprodine  | 9602 |
| G. Alphacetylmethadol (except<br>levoalphacetylmethadol<br>also known as levo-alpha-<br>acetylmethadol levothadyl  |      |

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**By JCAR at 10:53 am, Sep 25, 2023**

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|---|-------------|
| acetate or LAAM)  | 9603        |
| H. Alphameprodine   | 9604        |
| I. Alphamethadol  | 9605        |
| J. Alpha-methylfentanyl<br>(N-1-(alphamethyl-beta-phenyl) ethyl-4-piperidyl) propionanilide; 1-(1-methyl-2-phenylethyl)-4 ((N-propanilido) piperidine)          | 9814        |
| K. Alpha-methylthiofentanyl<br>(N-(1-methyl-2-(2-thienyl) ethyl-4-piperidiny)-N-phenylpropanamide)  | 9832        |
| L. Benzethidine   | 9606        |
| M. Betacetylmethadol  | 9607        |
| N. Beta-hydroxyfentanyl<br>(N-(1-(2-hydroxy-2-phenethyl)-4-piperidiny)-N-phenylpropanamide)   | 9830        |
| O. Beta-hydroxy-3-methylfentanyl (other name: N-(1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidiny)-N-phenylpropanamide)  | 9831        |
| P. N-[1-[2-hydroxy-2-(thiophen-2-yl) ethyl]piperidin-4-yl]-N-phenylpropionamide<br>(other names: beta-hydroxythiofentanyl)                                      | 9836        |
| Q. Betameprodine  | 9608        |
| R. Betamethadol   | 9609        |
| S. <i>beta</i> -Methyl fentanyl<br>(N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl) propionamide (Other name: $\beta$ -methyl fentanyl)                           | 9856        |
| T. <i>beta</i> '-Phenyl fentanyl<br>(N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide (other names: $\beta$ '-phenyl fentanyl; 3-phenylpropanoyl fentanyl) | 9842        |
| U. Betaprodine  | 9611        |
| <b>V. brorophine (1-(1-(1-(4-bromophenyl) ethyl)piperidin-4-yl)-1,3-dihydro-2H-benzo[d]imidazol-2-one)</b>  | <b>9098</b> |
| <b>W. Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide)</b>   | <b>9822</b> |
| [V]X. Clonitazene   | 9612        |
| [W]Y. Crotonyl fentanyl ((E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide)   | 9844        |
| [X]Z. N-(1-phenethylpiperidin-4-yl)-N-Phenylcyclopentanecarboxamide<br>(other name: cyclopentyl   |             |

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| fentanyl)  | 9847 |
| [Y]AA. Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide)  | 9845 |
| [Z]BB. Dextromoramide  | 9613 |
| [AA]CC. Diampromide  | 9615 |
| [BB]DD. Diethylthiambutene   | 9616 |
| [CC]EE. Difenoxyin   | 9168 |
| [DD]FF. Dimenoxadol  | 9617 |
| [EE]GG. Dimepseptanol  | 9618 |
| [FF]HH. Dimethylthiambutene  | 9619 |
| [GG]II. Dioxaphetyl butyrate   | 9621 |
| [HH]JJ. Dipipanone   | 9622 |
| [II]KK. Ethylmethylthiambutene   | 9623 |
| [JJ]LL. Etonitazene  | 9624 |
| [KK]MM. Etoxadine  | 9625 |
| [LL]NN. Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-yl) (phenyl)carbamate)   | 9851 |
| [MM]OO. N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (other names: 4-fluoroisobutyryl fentanyl, para-fluoroisobutyryl fentanyl) | 9824 |
| [NN]PP. 2'-Fluoro ortho-fluorofentanyl (N-(1-(2-fluorophenethyl) piperidin-4-yl)-N-(2-fluorophenyl) propionamide (other names: 2'-fluoro 2-fluorofentanyl)   | 9855 |
| [OO]QQ. N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (other names: furanyl fentanyl)  | 9834 |
| [PP]RR. Furethidine  | 9626 |
| [QQ]SS. Hydroxypethidine   | 9627 |
| [RR]TT. N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (other name: isobutyryl fentanyl)  | 9827 |
| [SS]UU. Isotonitazene (N,N-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl) ethan-1-amine)  | 9614 |
| [TT]VV. Ketobemidone   | 9628 |
| [UU]WW. Levomoramide   | 9629 |
| [VV]XX. Levophenacymorphan   | 9631 |
| [WW]YY. Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide)   | 9825 |
| [XX]ZZ. 4'-Methyl acetyl fentanyl  |      |

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| (N-(1-(4-methylphenethyl) piperidin-4-yl)-N-phenylacetamide)   | 9819   |
| [YY]AAA. 3-Methylfentanyl (N-(3-methyl-1-(2-phenylethyl)-4-piperidyl)-N-phenylpropanamide), its optical and geometric isomers, salts, and salts of isomers                           | 9813   |
| [ZZ]BBB. 3-Methylthiofentanyl (N-(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl)-N-phenylpropanamide)  | 9833   |
| CCC. Metonitazene (N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine)  | 9757   |
| [AAA]DDD. Morpheridine   | 9632   |
| [BBB]EEE. MPPP (1-methyl-4-phenyl-4-propionoxypiperidine)  | 9661   |
| [CCC]FFF. MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl) piperazine)  | (9560) |
| [DDD]GGG. Noracymethadol   | 9633   |
| [EEE]HHH. Norlevorphanol   | 9634   |
| [FFF]III. Normethadone   | 9635   |
| [GGG]JJJ. Norpipanone  | 9636   |
| [HHH]KKK. N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (other name: ocfentanil) | 9838   |
| [III]LLL. <i>ortho</i> -Fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl) acrylamide)   | 9852   |
| [JJJ]MMM. <i>ortho</i> -Fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl) butyramide (other name: 2-fluorobutyryl fentanyl)                                   | 9846   |
| [KKK]NNN. <i>ortho</i> -Fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl) propionamide); other name: 2-fluorofentanyl)  | 9816   |
| [LLL]OOO. <i>ortho</i> -Fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide)  | 9853   |
| [MMM]PPP. <i>ortho</i> -Methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl) acetamide (other name:   |        |

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| 2-methyl acetylfentanyl)   | 9848 |
| [ <b>NNN</b> ] <b>QQQ</b> . <i>ortho</i> -Methyl methoxyacetyl fentanyl (2-methoxy- <i>N</i> -(2-methylphenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)acetamide (other name: 2-methyl methoxyacetyl fentanyl) | 9820 |
| [ <b>OOO</b> ] <b>RRR</b> . <i>N</i> -(4-chlorophenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)isobutyramide (other name: <i>para</i> -chloroisobutyryl fentanyl)  | 9826 |
| [ <b>PPP</b> ] <b>SSS</b> . <i>para</i> -Fluorobutyryl fentanyl ( <i>N</i> -(4-fluorophenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)butyramide)   | 9823 |
| [ <b>QQQ</b> ] <b>TTT</b> . <i>para</i> -fluorofentanyl( <i>N</i> -(4-fluorophenyl)- <i>N</i> -(1-(2-phenethyl)-4-piperidinyl)propanamide  | 9812 |
| [ <b>RRR</b> ] <b>UUU</b> . <i>para</i> -Fluoro furanyl fentanyl ( <i>N</i> -(4-fluorophenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)furan-2-carboxamide)   | 9854 |
| [ <b>SSS</b> ] <b>VVV</b> . <i>para</i> -Methoxybutyryl fentanyl ( <i>N</i> -(4-methoxyphenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)butyramide)   | 9837 |
| [ <b>TTT</b> ] <b>WWW</b> . <i>para</i> -Methylfentanyl ( <i>N</i> -(4-methylphenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)propionamide (other name: 4-methylfentanyl)                                       | 9817 |
| [ <b>UUU</b> ] <b>XXX</b> . PEPAP (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine)   | 9663 |
| [ <b>VVV</b> ] <b>YYY</b> . Phenadoxone  | 9637 |
| [ <b>WWW</b> ] <b>ZZZ</b> . Phenampromide  | 9638 |
| [ <b>XXX</b> ] <b>AAAA</b> . Phenomorphan  | 9647 |
| [ <b>YYY</b> ] <b>BBBB</b> . Phenoperidine   | 9641 |
| [ <b>ZZZ</b> ] <b>CCCC</b> . Phenyl fentanyl ( <i>N</i> -(1-phenethylpiperidin-4-yl)- <i>N</i> -phenylbenzamide (other name: benzoyl fentanyl)   | 9841 |
| [ <b>AAAA</b> ] <b>DDDD</b> . Piritramide  | 9642 |
| [ <b>BBBB</b> ] <b>EEEE</b> . Proheptazine   | 9643 |
| [ <b>CCCC</b> ] <b>FFFF</b> . Properidine  | 9644 |
| [ <b>DDDD</b> ] <b>GGGG</b> . Propiram   | 9649 |
| [ <b>EEEE</b> ] <b>HHHH</b> . Racemoramide   | 9645 |
| [ <b>FFFF</b> ] <b>IIII</b> . <i>N</i> -(1-phenethylpiperidin-4-yl)- <i>N</i> -phenyltetrahydrofuran-  |      |

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| 2-carboxamide, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers (other name: tetrahydrofuranyl fentanyl)                             | 9843 |
| <b>[GGGG]JJJJ. Thiofentanyl (N-phenyl-N-(1-(2-thienyl)ethyl-4-piperidinyl)-propanamide</b>   | 9835 |
| <b>[HHHH]KKKK. Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide (other names: 2-thiofuranyl fentanyl; thiophene fentanyl)</b> | 9839 |
| <b>[IIII]LLLL. Tilidine</b>  | 9750 |
| <b>[JJJJ]MMMM. Trimeperidine</b>   | 9646 |
| <b>NNNN. U-47700 (3,4-Dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide)</b>  | 9547 |
| <b>OOOO. N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (other name: valeryl fentanyl)</b>  | 9840 |
| <b>PPPP. Zipeprol (1-methoxy-3-[4-(2-methoxy-2-phenylethyl)piperazin-1-yl]-1-phenylpropan-2-ol)</b>  | 9873 |

2. Opium derivatives. 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:

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| A. Acetorphine                           | 9319 |
| B. Acetyldihydrocodeine                  | 9051 |
| C. Benzylmorphine                        | 9052 |
| D. Codeine methylbromide                 | 9070 |
| E. Codeine-N-Oxide                       | 9053 |
| F. Cyprenorphine                         | 9054 |
| G. Desomorphine                          | 9055 |
| H. Dihydromorphine                       | 9145 |
| I. Drotebanol                            | 9335 |
| J. Etorphine (except hydrochloride salt) | 9056 |
| K. Heroin                                | 9200 |
| L. Hydromorphanol                        | 9301 |
| M. Methyldesorphine                      | 9302 |
| N. Methyldihydromorphine                 | 9304 |
| O. Morphine methylbromide                | 9305 |
| P. Morphine methylsulfonate              | 9306 |
| Q. Morphine-N-Oxide                      | 9307 |
| R. Myrophine                             | 9308 |
| S. Nicocodeine                           | 9309 |
| T. Nicomorphine                          | 9312 |
| U. Normorphine                           | 9313 |
| V. Pholcodine                            | 9314 |
| W. Thebacon                              | 9315 |

3[. *Opiate Similar Synthetic Substances. Substances scheduled by the United States Drug Enforcement Administration as substances that share a pharmacological profile similar to fentanyl, morphine, and*

other synthetic opioids, unless specifically excepted or unless listed in another schedule. These substances are—

|   |      |
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| A. Butyryl fentanyl ( <i>N</i> -<br>(1-phenethylpiperidin-4-yl)-<br><i>N</i> -phenylbutyramide)             | 9822 |
| B. U-47700 (3,4-Dichloro-<br><i>N</i> -[2-(dimethylamino)<br>cyclohexyl]- <i>N</i> -<br>methylbenzamide)    | 9547 |
| C. <i>N</i> -(1-phenethylpiperidin-<br>4-yl)- <i>N</i> -phenylpentanamide<br>(other name: valeryl fentanyl) | 9840 |

4]. 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 substances 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 paragraph (1)(A)4. of this rule only, the term isomer includes the optical, position, and geometric isomers.):

|  |      |
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| A. Alpha-ethyltryptamine   | 7249 |
| Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethenamine; 3-(2-aminobutyl)indole; alpha-ET; and AET; |      |
| B. 4-bromo-2,5-dimethoxyamphetamine  | 7391 |
| Some trade or other names: 4-bromo-2, 5- dimethoxy-a-methylphenethylamine; 4-bromo- 2, 5-DMA;                                  |      |
| C. 4-bromo-2,5-dimethoxyphenethylamine   | 7392 |
| D. 2,5-dimethoxyamphetamine  | 7396 |
| Some trade or other names: 2,5-dimethoxy-amethylphenethylamine; 2,5-DMA;   |      |
| E. 2,5-dimethoxy-4-ethylamphetamine  | 7399 |
| Some trade or other names: DOET;   |      |
| F. 2,5-dimethoxy-4-(n)-propylthiophenethylamine<br>(other name: 2C-T-7)  | 7348 |
| G. 2-(2,5-Dimethoxy-4-(n)-propylphenyl)<br>ethanamine (2C-P)   | 7524 |
| H. 2-(2,5-Dimethoxy-4-ethylphenyl)<br>ethanamine (2C-E)  | 7509 |
| I. 2-(2,5-Dimethoxy-4-methylphenyl)<br>ethanamine (2C-D)   | 7508 |
| J. 2-(2,5-Dimethoxy-4-nitro-<br>phenyl) ethanamine (2C-N)  | 7521 |
| K. 2-(2,5-Dimethoxyphenyl)<br>ethanamine (2C-H)  | 7517 |
| L. 2-(4-Chloro-2,5-dimethoxyphenyl)<br>ethanamine (2C-C)   | 7519 |
| M. 2-(4-Ethylthio-2,5-dimethoxyphenyl)<br>ethanamine (2C-T-2)  | 7385 |
| N. 2-(4-Iodo-2,5-dimethoxyphenyl)<br>ethanamine (2C-I)   | 7518 |
| O. 2-(4-Isopropylthio)-2,5-dimethoxyphenyl)<br>ethanamine (2C-T-4)   | 7532 |
| P. 4-methoxyamphetamine  | 7411 |
| Some trade or other names: 4-methoxy-amethylphenethylamine; paramethoxyamphetamine; PMA;                                       |      |
| Q. 5-methoxy-3,4-methylenedioxyamphetamine   | 7401 |
| R. 4-methyl-2,5-dimethoxyamphetamine   | 7395 |
| Some trade and other names: 4-methyl-2, 5- dimethoxy-a-methylphenethylamine; DOM; and STP;                                     |      |
| S. 3,4-methylenedioxyamphetamine   | 7400 |
| T. 3,4-methylenedioxymetham-<br>phetamine(MDMA)  | 7405 |

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| U. 3,4-methylenedioxy- <i>N</i> -ethylamphetamine (also known as <i>N</i> -ethylalpha-methyl-3,4 (methylenedioxy) phenethylamine, <i>N</i> -ethyl MDA, MDE, and MDEA)  | 7404 |
| V. <i>N</i> -hydroxy-3,4-methylenedioxyamphetamine (also known as <i>N</i> -hydroxy-alpha-methyl-3,4 (methylenedioxy) phenethylamine and <i>N</i> -hydroxy MDA)  | 7402 |
| W. 3,4,5-trimethoxyamphetamine   | 7390 |
| X. 5-MeO-DMT or 5-methoxy- <i>N,N</i> -dimethyltryptamine  | 7431 |
| Y. Alpha-methyltryptamine  | 7432 |
| Z. Bufotenine  | 7433 |
| Some trade and other names: 3-( <i>b</i> -Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; <i>N,N</i> -dimethylserotonin; 5-hydroxy- <i>N,N</i> -dimethyltryptamine; mappine;  |      |
| AA. Diethyltryptamine  | 7434 |
| Some trade and other names: <i>N,N</i> -Diethyltryptamine; DET;  |      |
| BB. Dimethyltryptamine   | 7435 |
| Some trade or other names: DMT;  |      |
| CC. 5-methoxy- <i>N,N</i> -diisopropyltryptamine (other name: 5-MeODIPT)   | 7439 |
| DD. 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- <i>b</i> ] indole; Tabernanthe iboga;   |      |
| EE. Lysergic acid diethylamide   | 7315 |
| FF. Marihuana  | 7360 |
| Some trade or other names: marijuana;  |      |
| GG. Mescaline  | 7381 |
| HH. Parahexyl  | 7374 |
| Some trade or other names: 3-Hexyl-1- hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl- 6H-dibenzo[ <i>b,d</i> ]pyran; Synhexyl;  |      |
| II. Peyote   | 7415 |
| Meaning all parts of the plant presently classified botanically as <i>Lophophora williamsii</i> Lemaire, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salt, derivative, mixture, or preparation of such plant, its seeds, or extracts;   |      |
| JJ. <i>N</i> -ethyl-3-piperidyl benzilate  | 7482 |
| KK. <i>N</i> -methyl-3-piperidyl benzilate   | 7484 |
| LL. Psilocybin   | 7437 |
| MM. Psilocyn   | 7438 |
| NN. Tetrahydrocannabinols naturally contained in a plant of the genus <i>Cannabis</i> ( <i>cannabis</i> 7370 plant), as well as synthetic equivalents of the substances contained in the <i>cannabis</i> plant or in the resinous extractives of such plant, and/or synthetic substances, derivatives, and their isomers, or both, with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following: |      |
| (I) 1 <i>cis</i> or <i>trans</i> tetrahydrocannabinol and their optical isomers;   |      |
| (II) 6 <i>cis</i> or <i>trans</i> tetrahydrocannabinol and their optical isomers;  |      |
| (III) 3,4 <i>cis</i> or <i>trans</i> tetrahydrocannabinol and its optical isomers; and   |      |
| (IV) Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions are covered;   |      |
| OO. Ethylamine analog of phencyclidine   | 7455 |
| Some trade or other names: <i>N</i> -ethyl-1- phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, <i>N</i> -(1-phenylcyclohexyl)-ethylamine, cyclohexamine, PCE;   |      |



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| PP. Pyrrolidine analog of phencyclidine   | 7458 |
| Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine PCPy, PHP;  |      |
| QQ. Thiophene analog of phencyclidine   | 7470 |
| Some trade or other names: 1-(1-(2-thienyl)-cyclohexyl)-piperidine, 2-thienyl analog of phencyclidine, TPCP, TCP;   |      |
| RR. 1-(1-(2-thienyl)cyclohexyl) pyrrolidine   | 7473 |
| Some other names: TCPy;   |      |
| SS. Salvia divinorum  |      |
| TT. Salvinorin A  |      |
| UU. 3-Fluoromethcathinone   | 1233 |
| VV. 4-Fluoromethcathinone   | 1238 |
| WW. Mephedrone, or 4-methylmethcathinone  | 1248 |
| XX. Methylenedioxy-pyrovalerone, MDPV, or (1-(1,3-Benzodioxol-5-yl)-2-(1-pyrrolidinyl)-1-pentanone  | 7535 |
| YY. Methylone, or 3,4-Methylenedioxy-methcathinone  | 7540 |
| ZZ. Quinolin-8-yl 1-pentyl-1Hindole-3-carboxylate (PB-22; QUPIC)  | 7222 |
| AAA. Quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22)  | 7225 |
| BBB. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1Hindazole-3-carboxamide (AB-FUBINACA)   | 7012 |
| CCC. N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (ADB-PINACA)   | 7035 |
| DDD. (1-pentyl-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (other names: UR-144, 1-pentyl-3-(2,2,3,3-tetramethylcyclopropyl)indole)   | 7144 |
| EEE. [1-(5-fluoro-pentyl)-1Hindol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone (other names: 5-fluoro-UR-144, 5-F-UR-144, XLR11, 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropyl)indole) | 7011 |
| FFF. N-(1-adamantyl)-1-pentyl-1Hindazole-3-carboxamide (other names: APINACA, AKB48)  | 7048 |
| GGG. 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine   |      |

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|      | (other names: 251-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5)  | 7538 |      |
| HHH. | 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine<br>(other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82)   | 7537 |      |
| III. | 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine<br>(other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36)  | 7536 |      |
| JJJ. | 4-methyl-N-ethylcathinone<br>(other names: 4-MEC; 2-(ethylamino)-1-(4-methylphenyl)propan-1-one)  | 1249 |      |
| KKK. | 4-methyl- $\alpha$ -pyrrolidinopropiophenone,<br>(other names: 4-MePPP; MePPP; 4-methyl- $\alpha$ -pyrrolidinopropiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)-propan-1-one) | 7498 |      |
| LLL. | $\alpha$ -pyrrolidinopentio-phenone<br>(other names: $\alpha$ -PVP; $\alpha$ -pyrrolidinovalerophenone; 1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one)                                 |      | 7545 |
| MMM. | Butylone<br>(other names: bk-MBDB; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one)   | 7541 |      |
| NNN. | Pentedrone<br>(other names: $\alpha$ -methylaminovalerophenone; 2-(methylamino)-1-phenylpentan-1-one)   | 1246 |      |
| OOO. | Pentylone<br>(other names: bk-MBDP; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one)   | 7542 |      |
| PPP. | Naphyrone<br>(other names: naphthylpyrovalerone; 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one)   | 1258 |      |
| QQQ. | $\alpha$ -pyrrolidinobutio-phenone<br>(other names: $\alpha$ -PBP; 1-phenyl-2-(pyrrolidin-1-yl)butan-1-one)   | 7546 |      |
| RRR. | N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide<br>(other names: AB-CHMINACA)   | 7031 |      |
| SSS. | N-(1-amino-3-methyl-1-  |      |      |

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| oxobutan-2-yl)-1-pentyl-<br>1H-indazole-3-carboxamide<br>(other names:<br>AB-PINACA)   | 7023   |
| TTT. [1-(5-fluoropentyl)-<br>1H-indazol-3-yl](naphthalen-<br>1-yl)methanone<br>(other names:<br>THJ-2201)  | 7024   |
| UUU. N-(1-amino-3,3-dimethyl-<br>1-oxobutan-2-yl)-1-<br>(cyclohexylmethyl)-<br>1H-indazole-3-carboxamide<br>(other names: MAB-<br>CHMINACA;<br>ADB-CHMINACA) | 7032   |
| VVV. methyl 2-(1-(5-fluoropentyl)-<br>1H-indazole-3-carboxamido)-3,3-<br>dimethylbutanoate<br>(other names:<br>5F-ADB; 5F-MDMB-PINACA)                       | 7034   |
| WWW. methyl 2-(1-(5-fluoropentyl)-<br>1H-indazole-3-carboxamido)-3-<br>methylbutanoate<br>(other names: 5F-AMB)  | 7033   |
| XXX. N-(adamantan-1-yl)-1-(5-<br>fluoropentyl)-1H-indazole-<br>3-carboxamide<br>(other names: 5F-APINACA, 5F-<br>AKB48)                                      | 7049   |
| YYY. N-(1-amino-3,3-dimethyl-<br>1-oxobutan-2-yl)-1-(4-<br>fluorobenzyl)-1H-indazole-<br>3-carboxamide<br>(other names: ADB-FUBINACA)                        | 7010   |
| ZZZ. methyl 2-(1-(cyclohexylmethyl)-<br>1H-indole-3-carboxamido)-3,3-<br>dimethylbutanoate<br>(other names: MDMB-CHMICA,<br>MMB-CHMINACA)                    | 7042   |
| AAAA. methyl 2-(1-(4-fluorobenzyl)-<br>1H-indazole-3-carboxamido)-3,3-<br>dimethylbutanoate<br>(other names: MDMB-FUBINACA)                                  | 7020   |
| BBBB. methyl 2-(1-(4-fluorobenzyl)-1H-<br>indazole-3-carboxamido)-3-<br>methylbutanoate<br>(other names: FUB-AMB, MMB-<br>FUBINACA, AMB-FUBINACA)            | (7021) |
| CCCC. 1-(1,3-benzodioxol-<br>5-yl)-2-(ethylamino)<br>propan-1-one (ethylone)   | 7547   |
| DDDD. Naphthalen-1-yl 1-<br>(5-fluoropentyl)-1H-<br>indole-3-carboxylate<br>(other names:<br>NM2201; CBL2201)  | 7221   |

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| EEEE. N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide<br>(other name: 5F-AB-PINACA)   | 7025 |
| FFFF. 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide<br>(other names: 4-CN-CUMYLBUTINACA; 4-cyano-CUMYLBUTINACA; 4-CN-CUMYLBINACA; CUMYL-4CNBINACA; SGT-78) | 7089 |
| GGGG. methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3-methylbutanoate<br>(other names: MMB-CHMICA; AMB-CHMICA)  | 7044 |
| HHHH. 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridine-3-carboxamide<br>(other name: 5F-CUMYL-P7AICA)   | 7085 |
| IIII. N-ethylpentylone (other names: ephylone, 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one)  | 7543 |
| JJJJ. methyl 2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate<br>(4F-MDMB-BINACA, 4F-MDMB-BUTINACA)  | 7043 |
| KKKK. 1-(4-methoxyphenyl)-N-methylpropan-2-amine<br>(other names: <i>para</i> -methoxymethamphetamine, PMMA)  | 1245 |
| LLLL. ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate<br>(other name: 5F-EDMB-PINACA)  | 7036 |
| MMMM. methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate (other names: 5F-MDMB-PICA; 5F-MDMB-2201)   | 7041 |
| NNNN. N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-FLUOROBENZYL))   | 7047 |
| OOOO. 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (other names: 5F-CUMYL-PINACA; SGT-25)   | 7083 |
| PPPP. (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)  |      |

|       |   |             |      |
|-------|---|-------------|------|
|       | methanone (other name: FUB-144)   | 7014        |      |
| QQQQ. | <i>N</i> -Ethylhexedrone (other names:<br><i>α</i> -ethylaminohexanophenone; 2-<br>(ethylamino)-1-phenylhexan-1-one)  | 7246        |      |
| RRRR. | <i>α</i> -Pyrrolidinohexanophenone<br>(other names: <i>α</i> -PHP; <i>α</i> -<br>pyrrolidinohexanophenone;<br>1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one)  |             | 7544 |
| SSSS. | 4-Methyl- <i>α</i> -ethylaminopentiophenone<br>(other names: 4-MEAP; 2-(ethylamino)-1-<br>(4-methylphenyl)pentan-1-one)   | 7245        |      |
| TTTT. | 4'-Methyl- <i>α</i> -pyrrolidinohexiophenone<br>(other names: MPHP; 4'-methyl-<br><i>α</i> -pyrrolidinohexanophenone; 1-<br>(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-<br>1-one)                        | 7446        |      |
| UUUU. | <i>α</i> -Pyrrolidinoheptaphenone<br>(other names: PV8; 1-phenyl-2-<br>(pyrrolidin-1-yl)heptan-1-one)   | 7548        |      |
| VVVV. | 4'-Chloro- <i>α</i> -pyrrolidinovalerophenone<br>(other names: 4-chloro- <i>α</i> -PVP; 4'-chloro- <i>α</i> -<br>pyrrolidinopentiophenone; 1-(4-<br>chlorophenyl)-2-(pyrrolidin-1-yl) pentan-<br>1-one) | 7443        |      |
| WWWW. | 2-(ethylamino)-2-<br>(3-methoxyphenyl)cyclohexan-<br>1-one (methoxetamine, MXE)   | 7286        |      |
| XXXX. | <b>1-(1,3-benzodioxol-5-yl)-2-<br/>(ethylamino)butan-1-one<br/>(other names: eutylone;<br/>bk-EBDB)</b>   | <b>7549</b> |      |

[XXXX]YYYY. Synthetic cannabinoids: Unless specifically exempted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, or which contains their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

(I) Any compound structurally derived from 3-(1-naphthoyl)indole or 1Hindol-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, whether or not substituted in the naphthyl ring to any extent. Including, but not limited to:

- |   |      |
|---|------|
| (a) AM2201, or 1-(5-fluoropentyl)-3-<br>(1-naphthoyl)indole           | 7201 |
| (b) JWH-007, or 1-pentyl-2-methyl-<br>3-(1-naphthoyl)indole           |      |
| (c) JWH-015, or 1-propyl-2-methyl-3-(1-<br>naphthoyl)indole           |      |
| (d) JWH-018, or 1-pentyl-<br>3-(1-naphthoyl)indole                    | 7118 |
| (e) JWH-019, or 1-hexyl-<br>3-(1-naphthoyl)indole                     | 7019 |
| (f) JWH-073, or 1-butyl-<br>3-(1-naphthoyl)indole                     | 7173 |
| (g) JWH-081, or 1-pentyl-3-(4-methoxy-1-<br>naphthoyl)indole          | 7081 |
| (h) JWH-098, or 1-pentyl-2-methyl-3-(4-<br>methoxy-1-naphthoyl)indole |      |

- (i) JWH-122, or 1-pentyl-3-(4-methyl-1-naphthoyl)indole 7122
- (j) JWH-164, or 1-pentyl-3-(7-methoxy-1-naphthoyl)indole
- (k) JWH-200, or 1-(2-(4-(morpholinyl)ethyl))-3-(1-naphthoyl)indole 7200
- (l) JWH-210, or 1-pentyl-3-(4-ethyl-1-naphthoyl)indole
- (m) JWH-398, or 1-pentyl-3-(4-chloro-1-naphthoyl)indole 7398

(II) 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, whether or not substituted in the naphthyl ring to any extent;

(III) Any compound structurally derived from 1-(1-naphthylmethyl)indene by substitution at the 3-position of the indene 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 indene ring to any extent, whether or not substituted in the naphthyl ring to any extent;

(IV) 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, whether or not substituted in the phenyl ring to any extent. Including, but not limited to:

- (a) JWH-201, or 1-pentyl-3-(4-methoxyphenylacetyl)indole
- (b) JWH-203, or 1-pentyl-3-(2-chlorophenylacetyl)indole 7203
- (c) JWH-250, or 1-pentyl-3-(2-methoxyphenylacetyl)indole 6250
- (d) JWH-251, or 1-pentyl-3-(2-methylphenylacetyl)indole
- (e) RCS-8, or 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole 7008

(V) 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, but not limited to:

(a) CP 47,497 & homologues, or 2-[(1*R*,3*S*)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol, where side chain n=5, and homologues where side chain n=4, 6, or 7 7297, 7298

(VI) Any compound containing 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, but not limited to:

- (a) AM-694, or 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole 7694
- (b) RCS-4, or 1-pentyl-3-(4-methoxybenzoyl)indole (SR-19 and RCS-4) 7104

(VII) CP 50,556-1, or [(6*S*,6*aR*,9*R*,10*aR*)-9-hydroxy-6-methyl-3-[(2*R*)-5-phenylpentan-2-yl]oxy-5,6,6*a*,7,8,9,10,10*a*-octahydrophenanthridin-1-yl] acetate;

(VIII) HU-210, or (6*aR*,10*aR*)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6*a*,7,10,10*a*-tetrahydrobenzo[*c*]chromen-1-ol;

(IX) HU-211, or Dexanabinol, (6*aS*,10*aS*)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6*a*,7,10,10*a*-tetrahydrobenzo[*c*]chromen-1-ol;

(X) Dimethylheptylpyran, or DMHP.

[5]4. 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:

- A. Gamma-hydroxybutyric acid and other names GHB;  
gamma-hydroxybutyrate; 4-hydroxybutyrate;  
4-hydroxybutonic acid; sodium oxybate;  
sodium oxybutyrate 2010
- B. Mecloqualone 2572
- C. Methaqualone 2565

[6]5. 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, and salts of isomers:

- A. Amineptine (7-[(10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-yl)amino]heptanoic acid) 1219**

[A]B. Aminorex 1585

Some trade or other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine;

[B]C. *N*-benzylpiperazine (some other names: BZP, 1-benzylpiperzaine) 7493

[C]D. Cathinone (Some trade or other names: 2-amino-1-phenyl-1-propanone, alphaaminopropiophenone, 2-aminopropiophenone and norephedrone) 1235

[D]E. 4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-oxazolamine; 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine) 1595

[E]F. Fenethylline 1503

**G. Mesocarb (N-phenyl-N'-(3-(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-yl)carbamide) 1227**

[F]H. Methcathinone 1237

Some trade or other names: 2-(methylamino)-propiofenone; alpha-(methylamino) propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-*N*-methylaminopropiophenone; monomethylpropion; ephedrone; *N*-methylcathinone; methylcathinine; AL-464; AL-422; AL-463 and URI 432;

**I. Methiopropamine (N-methyl-1-(thiophen-2-yl)propan-2-amine) 1478**

[G]J. 4-methoxymethcathinone

[H]K. *cis*-4-methylaminorex (*cis*-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) 1590

[I]L. 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP

[J]M. *N*-ethylamphetamine 1475

[K]N. *N,N*-dimethylamphetamine 1480

(some other names: *N,N*-alpha-trimethylbenzeneethanamine; *N,N*-alpha-trimethylphenethylamine)

[7]6. A temporary listing of substances subject to emergency scheduling under federal law shall include any material, compound, mixture, or preparation which contains any quantity of the following substances:

A. Fentanyl-related substances,  
their isomers, esters, ethers,  
salts, and salts of isomers,  
esters, and ethers.

9850

(I) Fentanyl-related substance means any substance not otherwise listed under another Administration Controlled Substance Code Number, 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 (1) or more of the following modifications:

(a) Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;

(b) Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino, or nitro groups;

(c) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxy, ester, ether, hydroxyl, halo, haloalkyl, amino, or nitro groups;

(d) Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; and/or

(e) Replacement of the *N*-propionyl group by another acyl group.

B. [1-(1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2H-benzo[d]imidazol-2-one, its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers

(Other names: bromphine;

1-[1-[1-(4-bromophenyl)ethyl]-4-piperidinyl]-1,3-dihydro-2H-benzimidazol-2-one)

9098

C. ]2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)-*N*, *N*-diethylethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

(other name: Butonitazene)

9751

[D]C. 2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-*N*, *N*-diethylethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other names: Etodesnitazene; etazene)

9765

[E]D. *N*, *N*-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

(other name: Flunitazene)

9756

[F]E. *N*,*N*-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

(other name: Metodesnitazene)

9764

[G. *N*, *N*-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers

(other name: Metonitazene)

9757]

[H]F. 2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-benzimidazole, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (other names: *N*-pyrrolidino



|   |      |
|---|------|
| etonitazene; etonitazepyne)   | 9758 |
| [I]G. <i>N, N</i> -diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1 <i>H</i> -benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers<br>(other name: Protonitazene) | 9759 |
| H. 4-(2-chlorophenyl)-2-ethyl-9-methyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine, its salts, isomers, and salts of isomers<br>(Other name: etizolam)                        | 2780 |
| I. 8-chloro-6-(2-fluorophenyl)-1-methyl-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine, its salts, isomers, and salts of isomers<br>(Other name: flualprazolam)                       | 2785 |
| J. 6-(2-chlorophenyl)-1-methyl-8-nitro-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine, its salts, isomers, and salts of isomers<br>(Other name: clonazolam)                           | 2786 |
| K. 8-bromo-6-(2-fluorophenyl)-1-methyl-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine, its salts, isomers, and salts of isomers<br>(Other name: flubromazolam)                        | 2788 |
| L. 7-chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2 <i>H</i> benzo[ <i>e</i> ][1,4]diazepin-2-one, its salts, isomers, and salts of isomers<br>(Other name: diclazepam)   | 2789 |

[8]7. Khat, to include all parts of the plant presently classified botanically as *catha edulis*, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seed, or extracts. 7032

(B) 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.

1. Substances, vegetable origin, or chemical synthesis. Unless specifically excepted or unless listed in another schedule, Schedule II shall include 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 a combination of extraction and chemical synthesis:

A. Opium and opiate; and any salt, compound, derivative, or preparation of opium or opiate, excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, naloxegol, naloxone, and naltrexone and their respective salts, but including the following:

|                         |      |
|-------------------------|------|
| (I) Raw opium           | 9600 |
| (II) Opium extracts     | 9610 |
| (III) Opium fluid       | 9620 |
| (IV) Powdered opium     | 9639 |
| (V) Granulated opium    | 9640 |
| (VI) Tincture of opium  | 9630 |
| (VII) Codeine           | 9050 |
| (VIII) Dihydroetorphine | 9334 |

|                             |      |
|-----------------------------|------|
| (IX) Ethylmorphine          | 9190 |
| (X) Etorphine hydrochloride | 9059 |
| (XI) Hydrocodone            | 9193 |
| (XII) Hydromorphone         | 9150 |
| (XIII) Metopon              | 9260 |
| (XIV) Morphine              | 9300 |
| (XV) Oripavine              | 9330 |
| (XVI) Oxycodone             | 9143 |
| (XVII) Oxymorphone          | 9652 |
| (XVIII) Thebaine            | 9333 |

B. Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in subparagraph (1)(B)1.A. of this rule shall be included in Schedule II, except that these substances shall not include the isoquinoline alkaloids of opium;

C. Opium poppy and poppy straw 9650

D. Coca leaves (9040) and any salt, compound, derivative, or preparation of coca leaves (including cocaine (9041) and ecgonine (9180) and their salts, isomers, derivatives, and salts of isomers and derivatives), and any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of these substances, except that the substances shall not include:

(I) Decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine; or

(II) Ioflupane;

E. Concentrate of 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

2. Opiates. 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 designation, dextrorphan, and levopropoxyphene excepted:

|  |      |
|--|------|
| A. Alfentanil                                    | 9737 |
| B. Alphaprodine                                  | 9010 |
| C. Anileridine                                   | 9020 |
| D. Bezitramide                                   | 9800 |
| E. Bulk Dextropropoxyphene<br>(Non-dosage Forms) | 9273 |
| F. Carfentanil                                   | 9743 |
| G. Dihydrocodeine                                | 9120 |
| H. Diphenoxylate                                 | 9170 |
| I. Fentanyl                                      | 9801 |
| J. Isomethadone                                  | 9226 |
| K. Levo-alphaacetylmethadol                      |      |

Some other names: levo-alphaacetylmethadol, levomethadyl acetate, LAAM 9648

|   |      |
|---|------|
| L. Levomethorphan   | 9210 |
| M. Levorphanol  | 9220 |
| N. Metazocine   | 9240 |
| O. Methadone  | 9250 |
| P. Methadone-Intermediate,<br>4-cyano-2-dimethylamino-<br>4,4-diphenyl butane | 9254 |
| Q. Moramide-Intermediate, 2-<br>methyl-3-morpholino-1,                        |      |

|   |      |
|---|------|
| 1-diphenylpropane-carboxylic acid   | 9802 |
| R. Oliceridine (N-[(3-methoxythiophen-2-yl)methyl]({2-[(9R)-9-(pyridin-2-yl)-6-oxaspiro[4.5]decan-9-yl]ethyl})amine fumarate) | 9245 |
| S. Pethidine (Meperidine)   | 9230 |
| T. Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine  | 9232 |
| U. Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate   | 9233 |
| V. Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid  | 9234 |
| W. Phenazocine  | 9715 |
| X. Piminodine   | 9730 |
| Y. Racemethorphan   | 9732 |
| Z. Racemorphan  | 9733 |
| AA. Remifentanil  | 9739 |
| BB. Sufentanil  | 9740 |
| CC. Tapentadol  | 9780 |
| DD. Thiafentanil  | 9729 |

3. 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:

|  |      |
|--|------|
| A. Amphetamine, its salts, optical isomers, and salts of its optical isomers | 1100 |
| B. Lisdexamfetamine, its salts, isomers, and salts of its isomers            | 1205 |
| C. Methamphetamine, its salts, isomers, and salts of its isomers             | 1105 |
| D. Phenmetrazine and its salts   | 1631 |
| E. Methylphenidate   | 1724 |

4. 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:

|                  |      |
|------------------|------|
| A. Amobarbital   | 2125 |
| B. Glutethimide  | 2550 |
| C. Pentobarbital | 2270 |
| D. Phencyclidine | 7471 |
| E. Secobarbital  | 2315 |

5. Hallucinogenic substances:

|             |      |
|-------------|------|
| A. Nabilone | 7379 |
|-------------|------|

Another name for nabilone: ( $\pm$ )trans-3-(1, 1-dimethylheptyl)-6, 6a,7,8,10,10a-hexahydro- 1-hydroxy-6, 6-dimethyl-9H-dibenzo(b,d) pyran-9-one.

B. Dronabinol [(-)-delta-9-*trans* tetrahydrocannabinol] in an oral solution in a drug product approved for marketing by the United States Food and Drug Administration. (7365)

6. Immediate precursors. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

A. Immediate precursor to amphetamine and methamphetamine:

- (I) Phenylacetone 8501
- Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;
- B. Immediate precursors to phencyclidine (PCP):
  - (I) 1-phenylcyclohexylamine 7460
  - (II) 1-piperidinocyclohexanecarbonitrile (PCC) 8603
- C. Immediate precursor to fentanyl:
  - (I) 4-anilino-*N*-phenethyl-4-piperidine (ANPP) 8333
  - (II) *N*-phenyl-*N*-(piperidin-4-yl)propionamide (norfentanyl) 8366

7. Any material, compound, mixture, or preparation which contains any quantity of the following alkyl nitrites:

- A. Amyl nitrite;
- B. Butyl nitrite.

(C) 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.

1. 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:

A. Those compounds, mixtures, or preparations in dosage unit form containing any stimulant substances listed in Schedule II which compounds, mixtures, or preparations were listed on August 25, 1971, as excepted compounds under 21 CFR 308.32 and any other drug of the quantitative composition shown in that list for those drugs or which is the same except that it contains a lesser quantity of controlled substances 1405

- B. Benzphetamine 1228
- C. Chlorphentermine 1645
- D. Clortermine 1647
- E. Phendimetrazine 1615

2. 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:

A. Any compound, mixture, or preparation containing—

- (I) Amobarbital 2126
- (II) Secobarbital 2316
- (III) Pentobarbital 2271

or any salt thereof and one (1) or more other active medicinal ingredients which are not listed in any schedule;

B. 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;

C. Any substance which contains any quantity of a derivative of barbituric acid or any salt thereof

|  |      |
|--|------|
| D. Chlorhexadol  | 2510 |
| E. Embutramide   | 2020 |
| F. Any drug product containing gamma hydroxybutyric acid, including its salts, isomers, and salts of isomer, for which an application is approved under section 505 of the Federal Food, Drug, and Cosmetic Act; | 2012 |
| G. Ketamine, its salts, isomer, and salts of isomers (some other names for ketamine: (±)-2-(2-chlorophenyl)-2-(methylamino)-cyclohexanone)   | 7285 |
| H. Lysergic acid   | 7300 |
| I. Lysergic acid amide   | 7310 |
| J. Methyprylon   | 2575 |
| K. Perampanel, and its salts, isomers, and salts of isomers  | 2261 |
| L. Sulfondiethylmethane  | 2600 |
| M. Sulfonethylmethane  | 2605 |
| N. Sulfonmethane   | 2610 |
| O. Tiletamine and zolazepam or any salt thereof  | 7295 |

Some trade or other names for a tiletaminezolazepam 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(1H)-one, flupyrazapon.

3. Nalorphine 9400

4. Narcotics drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof:

A. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium 9803

B. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9804

C. Not more than one and eight tenths grams (1.8gm) of dihydrocodeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9807

D. Not more than three hundred milligrams (300 mg) of ethylmorphine per one hundred milliliters (100 mL) or not more than fifteen milligrams (15 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9808

E. Not more than five hundred milligrams (500 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm) or not more than twenty-five milligrams (25 mg) per dosage unit, with one (1) or more active nonnarcotic ingredients in recognized therapeutic amounts 9809

F. Not more than fifty milligrams (50 mg) of morphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm), with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9810

5. Any material, compound, mixture, or preparation containing any of the following narcotic drugs or their salts, as set forth below:

A. Buprenorphine 9064

6. Anabolic steroids. [Unless specially excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation. DEA has assigned code 4000 for all anabolic steroids. Anabolic steroids. Any drug or hormonal substance, chemically and pharmacologically related to testosterone (other than estrogens, progestins, corticosteroids, and dehydroepiandrosterone) that promotes muscle growth, except an anabolic steroid which is expressly intended for administration through implants to cattle or other nonhuman species and which has been approved by the Secretary of Health and Human Services for that administration. If any person prescribes, dispenses, or distributes such steroid for human use, such person shall be considered to have prescribed, dispensed, or distributed an anabolic steroid within the meaning of this paragraph. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, esters, and ethers:

- A.  $3\beta,17\beta$ -dihydroxy- $5\alpha$ -androstane
- B.  $3\alpha,17\beta$ -dihydroxy- $5\alpha$ -androstane
- C.  $5\alpha$ -androstan- $3,17$ -dione
- D. 1-androstenediol ( $3\beta,17\beta$ -dihydroxy- $5\alpha$ -androst- $1$ -ene)
- E. 1-androstenediol ( $3\alpha,17\beta$ -dihydroxy- $5\alpha$ -androst- $1$ -ene)
- F. 4-androstenediol ( $3\beta,17\beta$ -dihydroxy-androst- $4$ -ene)
- G. 5-androstenediol ( $3\beta,17\beta$ -dihydroxy-androst- $5$ -ene)
- H. 1-androstenedione ( $[5\alpha]$ -androst- $1$ -en- $3,17$ -dione)
- I. 4-androstenedione (androst- $4$ -en- $3,17$ -dione)
- J. 5-androstenedione (androst- $5$ -en- $3,17$ -dione)
- K. Bolasterone ( $7\alpha,17\alpha$ -dimethyl- $17\beta$ -hydroxyandrost- $4$ -en- $3$ -one)
- L. Boldenone ( $17\beta$ -hydroxyandrost- $1,4$ -diene- $3$ -one)
- M. Boldione (androstra- $1,4$ -diene- $3,17$ -dione)
- N. Calusterone ( $7\beta,17\alpha$ -dimethyl- $17\beta$ -hydroxyandrost- $4$ -en- $3$ -one)
- O. Clostebol ( $4$ -chloro- $17\beta$ -hydroxyandrost- $4$ -en- $3$ -one)
- P. Dehydrochloromethyltestosterone ( $4$ -chloro- $17\beta$ -hydroxy- $17\alpha$ -methyl-androst- $1,4$ -dien- $3$ -one)
- Q. Desoxymethyltestosterone ( $17\alpha$ -methyl- $5\alpha$ -androst- $2$ -en- $17\beta$ -ol) (a.k.a. madol)
- R.  $\Delta 1$ -dihydrotestosterone (a.k.a. '1-testosterone') ( $17\beta$ -hydroxy- $5\alpha$ -androst- $1$ -en- $3$ -one)
- S. 4-dihydrotestosterone ( $17\beta$ -hydroxy-androstan- $3$ -one)
- T. Drostanolone ( $17\beta$ -hydroxy- $2\alpha$ -methyl- $5\alpha$ -androstan- $3$ -one)
- U. Ethylestrenol ( $17\alpha$ -ethyl- $17\beta$ -hydroxyestr- $4$ -ene)
- V. Fluoxymesterone ( $9$ -fluoro- $17\alpha$ -methyl- $11\beta,17\beta$ -dihydroxyandrost- $4$ -en- $3$ -one)
- W. Formebolone (Formebolone) ( $2$ -formyl- $17\alpha$ -methyl- $11\alpha,17\beta$ -dihydroxyandrost- $1,4$ -dien- $3$ -one)
- X. Furazabol ( $17\alpha$ -methyl- $17\beta$ -hydroxyandrostano[ $2,3$ -c]-furazan)
- Y.  $13\beta$ -ethyl- $17\beta$ -hydroxygon- $4$ -en- $3$ -one
- Z. 4-hydroxytestosterone ( $4,17\beta$ -dihydroxy-androst- $4$ -en- $3$ -one)
- AA. 4-hydroxy- $19$ -nortestosterone ( $4,17\beta$ -dihydroxy-estr- $4$ -en- $3$ -one)
- BB. Mestanolone ( $17\alpha$ -methyl- $17\beta$ -hydroxy- $5\alpha$ -androstan- $3$ -one)
- CC. Mesterolone ( $1\alpha$ -methyl- $17\beta$ -hydroxy- $[5\alpha]$ -androstan- $3$ -one)
- DD. Methandienone ( $17\alpha$ -methyl- $17\beta$ -hydroxyandrost- $1,4$ -dien- $3$ -one)
- EE. Methandriol ( $17\alpha$ -methyl- $3\beta,17\beta$ -dihydroxyandrost- $5$ -ene)
- FF. Methasterone ( $2\alpha,17\alpha$ -dimethyl- $5\alpha$ -androstan- $17\beta$ -ol- $3$ -one)
- GG. Methenolone ( $1$ -methyl- $17\beta$ -hydroxy- $5\alpha$ -androst- $1$ -en- $3$ -one)
- HH.  $17\alpha$ -methyl- $3\beta,17\beta$ -dihydroxy- $5\alpha$ -androstane
- II.  $17\alpha$ -methyl- $3\alpha,17\beta$ -dihydroxy- $5\alpha$ -androstane
- JJ.  $17\alpha$ -methyl- $3\beta,17\beta$ -dihydroxyandrost- $4$ -ene
- KK.  $17\alpha$ -methyl- $4$ -hydroxynandrolone ( $17\alpha$ -methyl- $4$ -hydroxy- $17\beta$ -hydroxyestr- $4$ -en- $3$ -one)

LL. Methyldienolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestra-4,9(10)-dien-3-one)  
MM. Methyltrienolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestra-4,9,11-trien-3-one)  
NN. Methyltestosterone (17 $\alpha$ -methyl-17-hydroxyandrost-4-en-3-one)  
OO. Mibolerone (7 $\alpha$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxyestr-4-en-3-one)  
PP. 17 $\alpha$ -methyl- $\Delta$ 1-dihydrotestosterone (17 $\beta$ -hydroxy-17 $\alpha$ -methyl-5 $\alpha$ -androst-1-en-3-one) (a.k.a. 17- $\alpha$ -methyl-1-testosterone)  
QQ. Nandrolone (17 $\beta$ -hydroxyestr-4-ene-3-one)  
RR. 19-nor-4-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxyestr-4-ene)  
SS. 19-nor-4-andro stenediol (3 $\alpha$ ,17 $\beta$ -dihydroxyestr-4-ene)  
TT. 19-nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione)  
UU. 19-nor-5-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxyestr-5-ene)  
VV. 19-nor-5-androstenediol (3 $\alpha$ ,17 $\beta$ -dihydroxyestr-5-ene)  
WW. 19-nor-4-androstenedione (estr-4-en-3,17-dione)  
XX. 19-nor-5-androstenedione (estr-5-en-3,17-dione)  
YY. Norbolethone (13 $\beta$ ,17 $\alpha$ -diethyl-17 $\beta$ -hydroxygon-4-en-3-one)  
ZZ. Norclostebol (4-chloro-17 $\beta$ -hydroxyestr-4-en-3-one)  
AAA. Norethandrolone (17 $\alpha$ -ethyl-17 $\beta$ -hydroxyestr-4-en-3-one)  
BBB. Normethandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestr-4-en-3-one)  
CCC. Oxandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-2-oxa-[5 $\alpha$ ]-androstan-3-one)  
DDD. Oxymesterone (17 $\alpha$ -methyl-4,17 $\beta$ -dihydroxyandrost-4-en-3-one)  
EEE. Oxymetholone (17 $\alpha$ -methyl-2-hydroxymethylene-17 $\beta$ -hydroxy-[5 $\alpha$ ]-androstan-3-one)  
FFF.Prostanozol (17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one-2-pyrazole)  
GGG. Stanolone ( $\Delta$ 1-dihydrotestosterone (a.k.a. 1-testosterone)(17 $\beta$ -hydroxy-5 $\alpha$ -androst-1-en-3-one))  
HHH. Stanozolol (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-[5 $\alpha$ ]-androst-2-eno[3,2-c]-pyrazole)  
III. Stenbolone (17 $\beta$ -hydroxy-2-methyl-[5 $\alpha$ ]-androst-1-en-3-one)  
JJJ. Testolactone(13-hydroxy-3-oxo-13,17-secoandrosta-1,4-dien-17-oic acid lactone)  
KKK. Testosterone(17 $\beta$ -hydroxyandrost-4-en-3-one);  
LLL.Tetrahydrogestrinone (13 $\beta$ ,17 $\alpha$ -diethyl-17 $\beta$ -hydroxygon-4,9, 11-trien-3-one)  
MMM. Trenbolone (17 $\beta$ -hydroxyestr-4,9,11-trien-3-one)  
NNN. Any salt, ester, or isomer of a drug or substance described or listed in this subparagraph, if that salt, ester, or isomer promotes muscle growth except an anabolic steroid which is expressly intended for administration through implants to cattle or other nonhuman species and which has been approved by the Secretary of Health and Human Services for that administration.]

**Unless specifically excepted or unless listed in another schedule, any substance meeting the definition of anabolic steroid as set forth in 21 CFR 1300.01, of the August 1, 2023, Code of Federal Regulations, as published by the Office of the Federal Resister, National Archives and Records Administration, 700 Pennsylvania Avenue, Washington, DC 20408 which is hereby incorporated by reference and does not include later amendments or additions, including any material, compound, mixture or preparation containing any quantity of the following substances, including its salts, esters and ethers (4000):**

- A. 5 $\alpha$ -androstan-3,17-dione;**
- B. 5 $\alpha$ -androstan-3,6,17-trione;**
- C. 1-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androst-1-ene);**
- D. 1-androstenediol (3 $\alpha$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androst-1-ene);**
- E. 4-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxy-androst-4-ene);**
- F. 5-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxy-androst-5-ene);**
- G. 1-androstenedione (5 $\alpha$ -androst-1-en-3,17-dione);**
- H. 4-androstenedione (androst-4-en-3,17-dione);**
- I. 5-androstenedione (androst-5-en-3,17-dione);**

J. bolasterone (7 $\alpha$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxyandrost-4-en-3-one);  
K. boldenone (17 $\beta$ -hydroxyandrost-1,4-diene-3-one);  
L. boldione (androsta-1,4-diene-3,17-dione);  
M. 6-bromo-androsta-1,4-diene-3,17-dione;  
N. 6-bromo-androstan-3,17-dione;  
O. calusterone (7 $\beta$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxyandrost-4-en-3-one);  
P. 4-chloro-17 $\alpha$ -methyl-androsta-1,4-diene-3,17 $\beta$ -diol;  
Q. 4-chloro-17 $\alpha$ -methyl-androst-4-ene-3 $\beta$ ,17 $\beta$ -diol;  
R. 4-chloro-17 $\alpha$ -methyl-17 $\beta$ -hydroxy-androst-4-en-3-one;  
S. 4-chloro-17 $\alpha$ -methyl-17 $\beta$ -hydroxy-androst-4-ene-3,11-dione;  
T. clostebol (4-chloro-17 $\beta$ -hydroxyandrost-4-en-3-one);  
U. dehydrochloromethyltestosterone (4-chloro-17 $\beta$ -hydroxy-17 $\alpha$ -methyl-androst-1,4-dien-3-one);  
V. desoxymethyltestosterone (17 $\alpha$ -methyl-5 $\alpha$ -androst-2-en-17 $\beta$ -ol) (a.k.a. "madol");  
W. 4-dihydrotestosterone (17 $\beta$ -hydroxy-androstan-3-one);  
X.  $\Delta$ 1-dihydrotestosterone (a.k.a. "1-testosterone") (17 $\beta$ -hydroxy-5 $\alpha$ -androst-1-en-3-one);  
Y. 3 $\beta$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androstan-3-one;  
Z. 3 $\alpha$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androstan-3-one;  
AA. 2 $\alpha$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxy-5 $\beta$ -androstan-3-one;  
BB. drostanolone (17 $\beta$ -hydroxy-2 $\alpha$ -methyl-5 $\alpha$ -androstan-3-one);  
CC. 2 $\alpha$ ,3 $\alpha$ -epithio-17 $\alpha$ -methyl-5 $\alpha$ -androstan-17 $\beta$ -ol;  
DD. estra-4,9,11-triene-3,17-dione;  
EE. 13 $\beta$ -ethyl-17 $\beta$ -hydroxygon-4-en-3-one;  
FF. ethylestrenol (17 $\alpha$ -ethyl-17 $\beta$ -hydroxyestr-4-ene);  
GG. fluoxymesterone (9-fluoro-17 $\alpha$ -methyl-11 $\beta$ ,17 $\beta$ -dihydroxyandrost-4-en-3-one);  
HH. formebolone (2-formyl-17 $\alpha$ -methyl-11 $\alpha$ ,17 $\beta$ -dihydroxyandrost-1,4-dien-3-one);  
II. furazabol (17 $\alpha$ -methyl-17 $\beta$ -hydroxyandrostan-2,3-c]furazan);  
JJ. [3,2-c]furazan-5 $\alpha$ -androstan-17 $\beta$ -ol;  
KK. 18 $\alpha$ -homo-3-hydroxy-estra-2,5(10)-dien-17-one;  
LL. 4-hydroxy-19-nortestosterone (4,17 $\beta$ -dihydroxy-estr-4-en-3-one);  
MM. 4-hydroxy-androst-4-ene-3,17-dione;  
NN. 17 $\beta$ -hydroxy-androstan-2,3-d]isoxazole;  
OO. 17 $\beta$ -hydroxy-androstan-3,2-c]isoxazole;  
PP. 3 $\beta$ -hydroxy-estra-4,9,11-trien-17-one;  
QQ. 4-hydroxytestosterone (4,17 $\beta$ -dihydroxy-androst-4-en-3-one);  
RR. mestanolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one);  
SS. mesterolone (1 $\alpha$ -methyl-17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one);  
TT. methandienone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyandrost-1,4-dien-3-one);  
UU. methandriol (17 $\alpha$ -methyl-3 $\beta$ ,17 $\beta$ -dihydroxyandrost-5-ene);  
VV. methasterone (2 $\alpha$ ,17 $\alpha$ -dimethyl-5 $\alpha$ -androstan-17 $\beta$ -ol-3-one or 2 $\alpha$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one);  
WW. methenolone (1-methyl-17 $\beta$ -hydroxy-5 $\alpha$ -androst-1-en-3-one);  
XX. 17 $\alpha$ -methyl-androsta-1,4-diene-3,17 $\beta$ -diol;  
YY. 17 $\alpha$ -methyl-5 $\alpha$ -androstan-17 $\beta$ -ol;  
ZZ. 17 $\alpha$ -methyl-androstan-3-hydroxyimine-17 $\beta$ -ol;  
AAA. 6 $\alpha$ -methyl-androst-4-ene-3,17-dione;  
BBB. 17 $\alpha$ -methyl-androst-2-ene-3,17 $\beta$ -diol;  
CCC. 17 $\alpha$ -methyl-3 $\beta$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androstan-3-one;  
DDD. 17 $\alpha$ -methyl-3 $\alpha$ ,17 $\beta$ -dihydroxy-5 $\alpha$ -androstan-3-one;  
EEE. 17 $\alpha$ -methyl-3 $\beta$ ,17 $\beta$ -dihydroxyandrost-4-ene;



FFF. 17 $\alpha$ -methyl-4-hydroxynandrolone (17 $\alpha$ -methyl-4-hydroxy-17 $\beta$ -hydroxyestr-4-en-3-one);

GGG. methyldienolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestra-4,9(10)-dien-3-one);

HHH. 17 $\alpha$ -methyl- $\Delta$ 1-dihydrotestosterone (17 $\beta$ -hydroxy-17 $\alpha$ -methyl-5 $\alpha$ -androst-1-en-3-one) (a.k.a. "17- $\alpha$ -methyl-1-testosterone");

III. methyltestosterone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyandrost-4-en-3-one);

JJJ. methyltrienolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestra-4,9,11-trien-3-one);

KKK. mibolerone (7 $\alpha$ ,17 $\alpha$ -dimethyl-17 $\beta$ -hydroxyestr-4-en-3-one);

LLL. nandrolone (17 $\beta$ -hydroxyestr-4-en-3-one);

MMM. 19-nor-4-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxyestr-4-ene);

NNN. 19-nor-4-androstenediol (3 $\alpha$ ,17 $\beta$ -dihydroxyestr-4-ene);

OOO. 19-nor-5-androstenediol (3 $\beta$ ,17 $\beta$ -dihydroxyestr-5-ene);

PPP. 19-nor-5-androstenediol (3 $\alpha$ ,17 $\beta$ -dihydroxyestr-5-ene);

QQQ. 19-nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione);

RRR. 19-nor-4-androstenedione (estr-4-en-3,17-dione);

SSS. 19-nor-5-androstenedione (estr-5-en-3,17-dione);

TTT. norbolethone (13 $\beta$ ,17 $\alpha$ -diethyl-17 $\beta$ -hydroxygon-4-en-3-one);

UUU. norclostebol (4-chloro-17 $\beta$ -hydroxyestr-4-en-3-one);

VVV. norethandrolone (17 $\alpha$ -ethyl-17 $\beta$ -hydroxyestr-4-en-3-one);

WWW. normethandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestr-4-en-3-one);

XXX. oxandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-2-oxa-5 $\alpha$ -androstan-3-one);

YYY. oxymesterone (17 $\alpha$ -methyl-4,17 $\beta$ -dihydroxyandrost-4-en-3-one);

ZZZ. oxymetholone (17 $\alpha$ -methyl-2-hydroxymethylene-17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one);

AAAA. prostanazol (17 $\beta$ -hydroxy-5 $\alpha$ -androstan[3,2-c]pyrazole or [3,2-c]pyrazole-5 $\alpha$ -androstan-17 $\beta$ -ol);

BBBB. [3,2-c]pyrazole-androst-4-en-17 $\beta$ -ol;

CCCC. stanozolol (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-5 $\alpha$ -androst-2-eno[3,2-c]-pyrazole);

DDDD. stenbolone (17 $\beta$ -hydroxy-2-methyl-5 $\alpha$ -androst-1-en-3-one);

EEEE. testolactone (13-hydroxy-3-oxo-13,17-secoandrosta-1,4-dien-17-oic acid lactone);

FFFF. testosterone (17 $\beta$ -hydroxyandrost-4-en-3-one);

GGGG. tetrahydrogestrinone (13 $\beta$ ,17 $\alpha$ -diethyl-17 $\beta$ -hydroxygon-4,9,11-trien-3-one); and

HHHH. trenbolone (17 $\beta$ -hydroxyestr-4,9,11-trien-3-one).

7. Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a United States Food and Drug Administration approved drug product 7369  
(Some other names for dronabinol: (6aRtrans)-6a,7,8,10a-tetrahydro-6.6.9-trimethyl-3-pentyl-6H-dibenzo (b,d) pyran-1-ol, or (-) -delta-9-(trans)-tetrahydrocannabinol.)

(D) 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 subsection. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.

1. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof:

A. Not more than one milligram (1 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit 9167

B. Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane) 9278

C. 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol, its salts, optical and geometric isomers, and salts of these isomers (including tramadol) 9752

D. 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 (1) 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:

(I) Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

(II) Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm); or

(III) Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm).

2. Depressants. 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:

|   |      |
|---|------|
| A. Alfaxalone                           | 2731 |
| B. Alprazolam                           | 2882 |
| C. Barbitol                             | 2145 |
| D. Brexanolone                          | 2400 |
| E. Bromazepam                           | 2748 |
| F. Camazepam                            | 2749 |
| G. Carisoprodol                         | 8192 |
| H. Chloral betaine                      | 2460 |
| I. Chloral hydrate                      | 2465 |
| J. Chlordiazepoxide                     | 2744 |
| K. Clobazam                             | 2751 |
| L. Clonazepam                           | 2737 |
| M. Clorazepate                          | 2768 |
| N. Clotiazepam                          | 2752 |
| O. Cloxazolam                           | 2753 |
| P. Daridorexant                         | 2410 |
| Q. Delorazepam                          | 2754 |
| R. Diazepam                             | 2765 |
| S. Dichloralphenazone                   | 2467 |
| T. Estazolam                            | 2756 |
| U. Ethchlorvynol                        | 2540 |
| V. Ethinamate                           | 2545 |
| W. Ethyl loflazepate                    | 2758 |
| X. Fludiazepam                          | 2759 |
| Y. Flunitrazepam                        | 2763 |
| Z. Flurazepam                           | 2767 |
| AA. Fospropofol                         | 2138 |
| BB. Halazepam                           | 2762 |
| CC. Haloxazolam                         | 2771 |
| DD. Ketazolam                           | 2772 |
| EE. Lemborexant                         | 2245 |
| FF. Loprazolam                          | 2773 |
| GG. Lorazepam                           | 2885 |
| HH. Lormetazepam                        | 2774 |
| II. Mebutamate                          | 2800 |
| JJ. Medazepam                           | 2836 |
| KK. Meprobamate                         | 2820 |
| LL. Methohexital                        | 2264 |
| MM. Methylphenobarbital (Mephobarbital) | 2250 |
| NN. Midazolam                           | 2884 |
| OO. Nimetazepam                         | 2837 |
| PP. Nitrazepam                          | 2834 |
| QQ. Nordiazepam                         | 2838 |
| RR. Oxazepam                            | 2835 |
| SS. Oxazolam                            | 2839 |

|                   |      |
|-------------------|------|
| TT. Paraldehyde   | 2585 |
| UU. Petrichloral  | 2591 |
| VV. Phenobarbital | 2285 |
| WW. Pinazepam     | 2883 |
| XX. Prazepam      | 2764 |
| YY. Quazepam      | 2881 |
| ZZ. Remimazolam   | 2846 |
| AAA. Suvorexant   | 2223 |
| BBB. Temazepam    | 2925 |
| CCC. Tetrazepam   | 2886 |
| DDD. Triazolam    | 2887 |
| EEE. Zaleplon     | 2781 |
| FFF. Zolpidem     | 2783 |
| GGG. Zopiclone    | 2784 |

3. [ Fenfluramine. 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:

A. Fenfluramine 1670

4.] Lorcaserin. 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:

A. Lorcaserin 1625

[5]4. 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, and salts of isomers:

|   |      |
|---|------|
| A. Cathine ((+)-norpseudoephedrine)   | 1230 |
| B. Diethylpropion   | 1610 |
| C. Fencamfamin  | 1760 |
| D. Fenproporex  | 1575 |
| E. Mazindol   | 1605 |
| F. Mefenorex  | 1580 |
| G. Modafinil  | 1680 |
| H. Pemoline (including organometallic complexes and chelates thereof)                               | 1530 |
| I. Phentermine  | 1640 |
| J. Pipradrol  | 1750 |
| K. Serdexmethylphenidate  | 1729 |
| L. Sibutramine  | 1675 |
| M. Solriamfetol (2-amino-3-phenylpropyl carbamate; benzenepropanol, beta-amino-, carbamate (ester)) | 1650 |
| N. SPA (-)-1-dimethylamino-1,2-diphenylethane   | 1635 |

[6]5. Other substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts:

|  |      |
|--|------|
| A. Pentazocine   | 9709 |
| B. Butorphanol (including its optical isomers)   | 9720 |
| C. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-oxopropyl] [(1S)-1-(4-phenyl-1 H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and salts of isomers | 9725 |

[7]6. Ephedrine. Any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system including their salts, isomers, and salts of isomers:

A. Ephedrine or its salts, optical isomers, or salts of optical isomers as the only active medicinal ingredient or contains ephedrine or its salts, optical isomers, or salts of optical isomers and therapeutically insignificant quantities of another active medicinal ingredient.

(E) 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 subsection.

1. Narcotic drugs containing nonnarcotic active medicinal ingredients. Any 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 follows, which shall include one (1) 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:

A. Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

B. Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

C. Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

D. Not more than two and five-tenths milligrams (2.5 mg) of diphenoxylate and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit;

E. Not more than one hundred milligrams (100 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm); and

F. Not more than five-tenths milligram (0.5 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit.

2. Stimulants. 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 stimulant effect on the central nervous system including its salts, isomers, and salts of isomers:

A. Pyrovalerone 1485

3. Any compound, mixture, or preparation containing any detectable quantity of pseudoephedrine or its salts or optical isomers, or salts of optical isomers or any compound, mixture, or preparation containing any detectable quantity of ephedrine or its salts or optical isomers, or salts of optical isomers if the drug preparations are starch-based solid dose forms, if such preparations are sold over the counter without a prescription. The following drug preparations containing ephedrine and pseudoephedrine are not scheduled controlled substances:

A. Drug preparations in liquid form; and

B. Drug preparations that require a prescription in order to be dispensed.

4. 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:

A. Ezogabine [N-[2-amino-4(4-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester] 2779

B. Ganaxolone (3 $\alpha$ -hydroxy-3 $\beta$ -methyl-5 $\alpha$ -pregnan-20-one) 2401

C. Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide] 2746

D. Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid] 2782

E. Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (also referred to as BRV; UCB-34714; Briviact) 2710

F. Lasmiditan [2,4,6-trifluoro-N-(6-(1-methylpiperidine-4-carbonyl) pyridine-2-yl)-benzamide] 2790

G. Cenobamate [(1R)-1-(2-chlorophenyl)-2-(tetrazol-

2-yl)ethyl] carbamate; 2H-tetrazole-2-ethanol, alpha-(2-chlorophenyl)-, carbamate (ester), (alphaR)-; carbamic acid (R)-(+)-1-(2-chlorophenyl)-2-(2H-tetrazol-2-yl)ethyl ester) 2720

(2) Excluded Nonnarcotic Substances. The following nonnarcotic substances which, under the Federal Food, Drug, and Cosmetic Act (21 U.S.C. 301) and section 201(g)(1) of the federal Controlled Substances Act (21 U.S.C. 811(g)(1)), may be lawfully sold over the counter without a prescription, are excluded from all schedules pursuant to section 195.015(5), RSMo.

*AUTHORITY: section 195.015, RSMo Supp. 2022, and section 195.195, RSMo 2016.\* Material found in this rule previously filed as 19 CSR 30-1.010. Original rule filed April 14, 2000, effective Nov. 30, 2000. Amended: Filed Jan. 31, 2003, effective July 30, 2003. Amended: Filed Sept. 30, 2016, effective May 30, 2017. Emergency amendment filed Oct. 25, 2018, effective Nov. 4, 2018, expired May 2, 2019. Amended: Filed Oct. 25, 2018, effective April 30, 2019. Emergency amendment filed Oct. 30, 2020, effective Nov. 16, 2020, expired May 14, 2021. Amended: Filed Oct. 30, 2020, effective April 30, 2021. Emergency amendment filed Sept. 28, 2021, effective Oct. 13, 2021, expired April 10, 2022. Amended: Filed Sept. 28, 2021, effective March 30, 2022. Emergency amendment filed Sept. 12, 2022, effective Oct. 3, 2022, expired March 31, 2023. Amended: Filed Sept. 12, 2022, effective March 30, 2023. . **Amended: Filed: September 25, 2023.***

*\*Original authority: 195.015, RSMo 1971, amended 1989, 2014, 2020, and 195.195, RSMo 1957, amended 1971, 1989, 1993, 2014.*

*PUBLIC COST: This proposed amendment will not cost state agencies or political subdivisions more than five hundred dollars (\$500) in the aggregate.*

*PRIVATE COST: This proposed amendment will not cost private entities more than five hundred dollars (\$500) in the aggregate.*

*NOTICE TO SUBMIT COMMENTS: Anyone may file a statement in support or in opposition to this proposed amendment with E.J. Jackson, Missouri Department of Health and Senior Services, Bureau of Narcotics and Dangerous Drugs, PO Box 570, Jefferson City, MO 65102, or via email at BNDD@health.mo.gov. To be considered, comments must be received within thirty (30) days after publication of this notice in the **Missouri Register**. No public hearing is scheduled.*