Title 21 —Food and Drugs
Chapter II —Drug Enforcement Administration, Department of Justice
Part 1308 —Schedules of Controlled Substances
Schedules

Authority: 21 U.S.C. 811, 812, 871(b), 956(b), unless otherwise noted.
Source: 38 FR 8254, Mar. 30, 1973, unless otherwise noted. Redesignated at 38 FR 26609, Sept. 24, 1973.

§ 1308.11 Schedule I.

Link to an amendment published at 90 FR 57542, Dec. 11, 2025.

(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.*** 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 (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

(1) Acetyl-*alpha*-methylfentanyl (*N*-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-*N*-phenylacetamide)	9815
(2) Acetylmethadol	9601
(3) Acetyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacetamide)	9821
(4) Acryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacrylamide; also known as acryloylfentanyl)	9811
(5) AH-7921 (3,4-dichloro-*N*-[(1-dimethylamino)cyclohexylmethyl]benzamide)	9551
(6) Allylprodine	9602
(7) Alphacetylmethadol (except *levo*-alphacetylmethadol also known as *levo-alpha*-acetylmethadol, levomethadyl acetate, or LAAM)	9603
(8) Alphameprodine	9604
(9) Alphamethadol	9605
(10) *alpha'*-Methyl butyryl fentanyl (2-methyl-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbutanamide)	9864
(11) *alpha*-Methylfentanyl (*N*-[1-(*alpha*-methyl-*beta*-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-2-phenylethyl)-4-(*N*-propanilido)piperidine)	9814
(12) *alpha*-Methylthiofentanyl (*N*-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-*N*-phenylpropanamide)	9832
(13) Benzethidine	9606
(14) Betacetylmethadol	9607

(15) *beta*-Hydroxyfentanyl (*N*-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-*N*-phenylpropanamide)	9830
(16) *beta*-Hydroxy-3-methylfentanyl (*N*-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidinyl]-*N*-phenylpropanamide)	9831
(17) *beta*-Hydroxythiofentanyl (*N*-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-*N*-phenylpropionamide)	9836
(18) Betameprodine	9608
(19) Betamethadol	9609
(20) *beta*-Methyl fentanyl (*N*-phenyl-*N*-(1-(2-phenylpropyl)piperidin-4-yl)propionamide; also known as β-methyl fentanyl)	9856
(21) *beta*-methylacetyl fentanyl (*N*-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)acetamide)	9868
(22) *beta'*-Phenyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*,3-diphenylpropanamide; also known as β'-phenyl fentanyl; 3-phenylpropanoyl fentanyl)	9842
(23) Betaprodine	9611
(24) brorphine (1-(1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2*H*-benzo[*d*]imidazol-2-one)	9098
(25) Butonitazene (2-(2-(4-butoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N*,*N*-diethylethan-1-amine)	9751
(26) Butyryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbutyramide)	9822
(27) Clonitazene	9612
(28) Crotonyl fentanyl ((*E*)-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbut-2-enamide)	9844
(29) Cyclopentyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylcyclopentanecarboxamide)	9847
(30) Cyclopropyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylcyclopropanecarboxamide)	9845
(31) Dextromoramide	9613
(32) Diampromide	9615
(33) Diethylthiambutene	9616
(34) Difenoxin	9168
(35) Dimenoxadol	9617
(36) Dimepheptanol	9618
(37) 2',5'-Dimethoxyfentanyl (*N*-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-*N*-phenylpropionamide)	9861
(38) Dimethylthiambutene	9619
(39) Dioxaphetyl butyrate	9621
(40) Dipipanone	9622
(41) Ethylmethylthiambutene	9623
(42) 2-(2-(4-ethoxybenzyl)-1*H*-benzimidazol-1-yl)-*N*,*N*-diethylethan-1-amine (Other names: etodesnitazene; etazene)	9765
(43) Etonitazene	9624
(44) Etoxeridine	9625
(45) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-yl)(phenyl)carbamate)	9851
(46) Flunitazene (*N*,*N*-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine)	9756
(47) 4-Fluoroisobutyryl fentanyl	9824

(*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide; also known as *para*-fluoroisobutyryl fentanyl)	
(48) 2'-Fluoro *ortho*-fluorofentanyl (*N*-(1-(2-fluorophenethyl)piperidin-4-yl)-*N*-(2-fluorophenyl)propionamide; also known as 2'-fluoro 2-fluorofentanyl)	9855
(49) Furanyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylfuran-2-carboxamide)	9834
(50) 3-Furanyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylfuran-3-carboxamide)	9860
(51) Furethidine	9626
(52) Hydroxypethidine	9627
(53) Isobutyryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylisobutyramide)	9827
(54) Isotonitazene (*N,N*-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine)	9614
(55) Isovaleryl fentanyl (3-methyl-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbutanamide)	9862
(56) Ketobemidone	9628
(57) Levomoramide	9629
(58) Levophenacylmorphan	9631
(59) *meta*-Fluorofentanyl *(N*-(3-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide)	9857
(60) *meta*-fluorofuranyl fentanyl (*N*-(3-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide)	9871
(61) *meta*-Fluoroisobutyryl fentanyl (*N*-(3-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide)	9858
(62) Methoxyacetyl fentanyl (2-methoxy-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacetamide)	9825
(63) 4'-Methyl acetyl fentanyl (*N*-(1-(4-methylphenethyl)piperidin-4-yl)-*N*-phenylacetamide)	9819
(64) 2-Methyl AP-237 (1-(2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl)butan-1-one)	9664
(65) 3-Methylfentanyl (*N*-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide)	9813
(66) 3-Methylthiofentanyl (*N*-[3-methyl-1-(2-thienylethyl)-4-piperidinyl]-*N*-phenylpropanamide)	9833
(67) Metodesnitazene (*N,N*-diethyl-2-(2-(4-methoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine)	9764
(68) Metonitazene (*N,N*-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine)	9757
(69) Morpheridine	9632
(70) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine)	9661
(71) MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine)	9560
(72) *N*-Desethyl isotonitazene (*N*-ethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine)	9760
(73) Noracymethadol	9633
(74) Norlevorphanol	9634
(75) Normethadone	9635
(76) Norpipanone	9636
(77) *N*-Piperidinyl etonitazene (2-(4-ethoxybenzyl)-5-nitro-1-(2-(piperidin-1-yl)ethyl)-1*H*-benzimidazole (other names: etonitazepipne)	9761

(78) 2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (Other names: *N*-pyrrolidino etonitazene; etonitazepyne)	9758
(79) 2-(4-methoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (Other names: *N*-pyrrolidino metonitazene; metonitazepyne)	9762
(80) 5-nitro-2-(4-propoxybenzyl)-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (other names: *N*-pyrrolidino protonitazene; protonitazepyne)	9763
(81) Ocfentanil (*N*-(2-fluorophenyl)-2-methoxy-*N*-(1-phenethylpiperidin-4-yl)acetamide)	9838
(82) *ortho*-chlorofentanyl (*N*-(2-chlorophenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide	9828
(83) *ortho*-Fluoroacryl fentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)acrylamide)	9852
(84) *ortho*-Fluorobutyryl fentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)butyramide; also known as 2-fluorobutyryl fentanyl)	9846
(85) *ortho*-Fluorofentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide); also known as 2-fluorofentanyl)	9816
(86) *ortho*-Fluorofuranyl fentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide)	9863
(87) *ortho*-Fluoroisobutyryl fentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide)	9853
(88) *ortho*-Methyl acetylfentanyl (*N*-(2-methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)acetamide; also known as 2-methyl acetylfentanyl)	9848
(89) *ortho*-methylcyclopropyl fentanyl (*N*-(2-methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide)	9849
(90) *ortho*-Methyl methoxyacetyl fentanyl (2-methoxy-*N*-(2-methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)acetamide; also known as 2-methyl methoxyacetyl fentanyl)	9820
(91) *para*-Chloroisobutyryl fentanyl (*N*-(4-chlorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide)	9826
(92) *para*-chlorofentanyl (*N*-(4-chlorophenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide)	9818
(93) *para*-Fluorobutyryl fentanyl (*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)butyramide)	9823
(94) *para*-Fluorofentanyl (*N*-(4-fluorophenyl)-*N*-[1-(2-phenylethyl)-4-piperidinyl]propanamide)	9812
(95) *para*-Fluoro furanyl fentanyl (*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide)	9854
(96) *para*-fluoro valeryl fentanyl (*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)pentanamide)	9870
(97) *para*-Methoxybutyryl fentanyl (*N*-(4-methoxyphenyl)-*N*-(1-phenethylpiperidin-4-yl)butyramide)	9837
(98) *para*-Methoxyfuranyl fentanyl (*N*-(4-methoxyphenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide	9859
(99) *para*-Methylcyclopropyl fentanyl (*N*-(4-methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide)	9865
(100) *para*-Methylfentanyl (*N*-(4-methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide; also known as 4-methylfentanyl)	9817
(101) PEPAP (1-(2-phenylethyl)-4-phenyl-4-acetoxypiperidine)	9663
(102) Phenadoxone	9637

(103) Phenampromide	9638
(104) Phenomorphan	9647
(105) Phenoperidine	9641
(106) Phenyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbenzamide; also known as benzoyl fentanyl)	9841
(107) Piritramide	9642
(108) Proheptazine	9643
(109) Properidine	9644
(110) Propiram	9649
(111) *N,N*-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: protonitazene)	9759
(112) Racemoramide	9645
(113) Tetrahydrofuranyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenyltetrahydrofuran-2-carboxamide)	9843
(114) tetrahydrothiofuranyl fentanyl (also known as: tetrahydrothiophene fentanyl) (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenyltetrahydrothiophene-2-carboxamide)	9869
(115) Thiofentanyl (*N*-phenyl-*N*-[1-(2-thienyl)ethyl-4-piperidinyl]propanamide)	9835
(116) Thiofuranyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylthiophene-2-carboxamide; also known as 2-thiofuranyl fentanyl; thiophene fentanyl)	9839
(117) Tilidine	9750
(118) Trimeperidine	9646
(119) U-47700 (3,4-dichloro-*N*-[2-(dimethylamino)cyclohexyl]-*N*-methylbenzamide)	9547
(120) Valeryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylpentanamide)	9840
(121) Zipeprol (1-methoxy-3-[4-(2-methoxy-2-phenylethyl)piperazin-1-yl]-1-phenylpropan-2-ol)	9873

(c) ***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:

(1) Acetorphine	9319
(2) Acetyldihydrocodeine	9051
(3) Benzylmorphine	9052
(4) Codeine methylbromide	9070
(5) Codeine-N-Oxide	9053
(6) Cyprenorphine	9054
(7) Desomorphine	9055
(8) Dihydromorphine	9145
(9) Drotebanol	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

(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 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 this paragraph only, the term "isomer" includes the optical, position and geometric isomers):

(1) Alpha-ethyltryptamine	7249
Some trade or other names: etryptamine; Monase; α-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; α-ET; and AET.	
(2) 4-bromo-2,5-dimethoxy-amphetamine	7391
Some trade or other names: 4-bromo-2,5-dimethoxy-α-methylphenethylamine; 4-bromo-2,5-DMA	
(3) 4-Bromo-2,5-dimethoxyphenethylamine	7392
Some trade or other names: 2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl DOB; 2C-B, Nexus.	
(4) 2,5-dimethoxyamphetamine	7396
Some trade or other names: 2,5-dimethoxy-α-methylphenethylamine; 2,5-DMA	
(5) 2,5-dimethoxy-4-ethylamphet-amine	7399
Some trade or other names: DOET	
(6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7)	7348
(7) 4-methoxyamphetamine	7411
Some trade or other names: 4-methoxy-α-methylphenethylamine; paramethoxyamphetamine, PMA	
(8) 5-methoxy-3,4-methylenedioxy-amphetamine	7401
(9) 4-methyl-2,5-dimethoxy-amphetamine	7395

Some trade and other names: 4-methyl-2,5-dimethoxy-α-methylphenethylamine; “DOM”; and “STP”

(10) 3,4-methylenedioxy amphetamine 7400

(11) 3,4-methylenedioxymethamphetamine (MDMA) 7405

(12) 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4(methylenedioxy)-phenethylamine, N-ethyl MDA, MDE, MDEA 7404

(13) N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-alpha-methyl-3,4(methylenedioxy)-phenethylamine, and N-hydroxy MDA 7402

(14) 3,4,5-trimethoxy amphetamine 7390

(15) 5-methoxy-N,N-dimethyltryptamine Some trade or other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT 7431

(16) Alpha-methyltryptamine (other name: AMT) 7432

(17) Bufotenine 7433

Some trade and other names: 3-(β-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine

(18) Diethyltryptamine 7434

Some trade and other names: N,N-Diethyltryptamine; DET

(19) Dimethyltryptamine 7435

Some trade or other names: DMT

(20) 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT) 7439

(21) Ibogaine 7260

Some trade and other names: 7-Ethyl-6,6β,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) Marihuana 7360

(24) Mescaline 7381

(25) Parahexyl—7374; some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl.

(26) 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))

(27) N-ethyl-3-piperidyl benzilate 7482

(28) N-methyl-3-piperidyl benzilate 7484

(29) Psilocybin 7437

(30) Psilocyn 7438

(31) Tetrahydrocannabinols 7370

(i) Meaning tetrahydrocannabinols, except as in paragraph (d)(31)(ii) of this section,

naturally contained in a plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of the substances contained in the cannabis plant, or in the resinous extractives of such plant, and/or synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following:

1 cis or trans tetrahydrocannabinol, and their optical isomers

6 cis or trans tetrahydrocannabinol, and their optical isomers

3, 4 cis or trans tetrahydrocannabinol, and its optical isomers

(Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions covered.)

(ii) Tetrahydrocannabinols does not include any material, compound, mixture, or preparation that falls within the definition of hemp set forth in 7 U.S.C. 1639*o*.

(32) Ethylamine analog of phencyclidine 7455

Some trade or other names: N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine, N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE

(33) Pyrrolidine analog of phencyclidine 7458

Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP

(34) Thiophene analog of phencyclidine 7470

Some trade or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP

(35) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine 7473

Some other names: TCPy

(36) 4-methylmethcathinone (Mephedrone) 1248

(37) 3,4-methylenedioxypyrovalerone (MDPV) 7535

(38) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E) 7509

(39) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D) 7508

(40) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C) 7519

(41) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I) 7518

(42) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2) 7385

(43) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4) 7532

(44) 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H) 7517

(45) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N) 7521

(46) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P) 7524

(47) 3,4-Methylenedioxy-N-methylcathinone (Methylone) 7540

(48) (1-pentyl-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (UR-144) (7144)

(49) [1-(5-fluoro-pentyl)-1*H*-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone (5-fluoro-UR-144, XLR11) (7011)

(50) *N*-(1-adamantyl)-1-pentyl-1*H*-indazole-3-carboxamide (APINACA, AKB48) (7048)

(51) quinolin-8-yl 1-pentyl-1*H*-indole-3-carboxylate (PB-22; QUPIC) (7222)

(52) quinolin-8-yl 1-(5-fluoropentyl)-1*H*-indole-3-carboxylate (5-fluoro-PB-22; 5F-PB-22) (7225)

(53) *N*-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (AB-FUBINACA) (7012)

(54) *N*-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1*H*-indazole-3-carboxamide (ADB-PINACA) (7035)

(55) 2-(4-iodo-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (25I-NBOMe, 2C-I-NBOMe) (7538)

(56) 2-(4-chloro-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (25C-NBOMe, 2C-C-NBOMe) (7537)

(57) 2-(4-bromo-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (25B-NBOMe, 2C-B-NBOMe) (7536)

(58) Marihuana Extract 7350

Meaning an extract containing one or more cannabinoids that has been derived from any plant of the genus Cannabis, containing greater than 0.3% delta-9-tetrahydrocannabinol on a dry weight basis, other than the separated resin (whether crude or purified) obtained from the plant.

(59) 4-methyl-*N*-ethylcathinone (4-MEC) (1249)

(60) 4-methyl-*alpha*-pyrrolidinopropiophenone (4-MePPP) (7498)

(61) *alpha*-pyrrolidinopentiophenone (α-PVP) (7545)

(62) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one (butylone, bk-MBDB) (7541)

(63) 2-(methylamino)-1-phenylpentan-1-one (pentedrone) (1246)

(64) 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one (pentylone, bk-MBDP) (7542)

(65) 4-fluoro-*N*-methylcathinone (4-FMC; flephedrone) (1238)

(66) 3-fluoro-*N*-methylcathinone (3-FMC) (1233)

(67) 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (naphyrone) (1258)

(68) *alpha*-pyrrolidinobutiophenone (α-PBP) (7546)

(69) *N*-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1*H*-indazole-3-carboxamide (AB-CHMINACA) (7031)

(70) *N*-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1*H*-indazole-3-carboxamide (AB-PINACA) (7023)

(71) [1-(5-fluoropentyl)-1*H*-indazol-3-yl](naphthalen-1-yl)methanone (THJ-2201) (7024)

(72) *N*-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1*H*-indazole-3-carboxamide (MAB-CHMINACA; ADB-CHMINACA) (7032)

(73) methyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5F-ADB; 5F-MDMB-PINACA) 7034

(74) methyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3-methylbutanoate (Other names: 5F-AMB) 7033

(75) *N*-(adamantan-1-yl)-1-(5-fluoropentyl)-1*H*-indazole-3-carboxamide (Other names: 5F-APINACA, 5F-AKB48) 7049

(76) *N*-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (Other names: ADB-FUBINACA) 7010

(77) methyl 2-(1-(cyclohexylmethyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (Other names: MDMB-CHMICA, MMB-CHMINACA) 7042

(78) methyl 2-(1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other names: MDMB-FUBINACA)	7020
(79) methyl 2-(1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamido)-3-methylbutanoate, (FUB-AMB, MMB-FUBINACA, AMB-FUBINACA)	(7021)
(80) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)propan-1-one (ethylone)	7547
(81) Naphthalen-1-yl 1-(5-fluoropentyl)-1*H*-indole-3-carboxylate (Other names: NM2201; CBL2201)	7221
(82) *N*-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1*H*-indazole-3-carboxamide (Other name: 5F-AB-PINACA)	7025
(83) 1-(4-cyanobutyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (Other names: 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINACA; 4-CN-CUMYL BINACA; CUMYL-4CN-BINACA; SGT-78)	7089
(84) methyl 2-(1-(cyclohexylmethyl)-1*H*-indole-3-carboxamido)-3-methylbutanoate (Other names: MMB-CHMICA; AMB-CHMICA)	7044
(85) 1-(5-fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-pyrrolo[2,3-b]pyridine-3-carboxamide (Other name: 5F-CUMYL-P7AICA)	7085
(86) *N*-ethylpentylone (Other names: ephylone, 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one)	7543
(87) methyl 2-(1-(4-fluorobutyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (4F-MDMB-BINACA, 4F-MDMB-BUTINACA)	7043
(88) 1-(4-methoxyphenyl)-*N*-methylpropan-2-amine (other names: *para*-methoxymethamphetamine, PMMA)	(1245)
(89) ethyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (other name: 5F-EDMB-PINACA)	7036
(90) methyl 2-(1-(5-fluoropentyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (other names: 5F-MDMB-PICA; 5F-MDMB-2201)	7041
(91) *N*-(adamantan-1-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (other names: FUB-AKB48; FUB-APINACA; AKB48 N-(4-FLUOROBENZYL))	7047
(92) 1-(5-fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (other names: 5F-CUMYL-PINACA; SGT-25)	7083
(93) (1-(4-fluorobenzyl)-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (other name: FUB-144)	7014
(94) *N*-Ethylhexedrone (Other names: α-ethylaminohexanophenone; 2-(ethylamino)-1-phenylhexan-1-one)	7246
(95) *alpha*-Pyrrolidinohexanophenone (Other names: α-PHP; α-pyrrolidinohexanophenone; 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one)	7544
(96) 4-Methyl-*alpha*-ethylaminopentiophenone (Other names: 4-MEAP; 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one)	7245
(97) 4'-Methyl-*alpha*-pyrrolidinohexiophenone (Other names: MPHP; 4'-methyl-*alpha*-pyrrolidinohexanophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one)	7446
(98) *alpha*-Pyrrolidinoheptaphenone (Other names: PV8; 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one)	7548
(99) 4'-Chloro-*alpha*-pyrrolidinovalerophenone (Other names: 4-chloro-α-PVP; 4'-chloro-α-pyrrolidinopentiophenone; 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one)	7443

(100) 2-(ethylamino)-2-(3-methoxyphenyl)cyclohexan-1-one (methoxetamine, MXE)	7286
(101) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)butan-1-one (other names: eutylone; bk-EBDB)	7549
(102) *N*-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1*H*-indazole-3-carboxamide (other name: ADB-BUTINACA)	7027
(103) 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one (other names: α-PiHP; *alpha*-PiHP)	7551
(104) 2-(methylamino)-1-(3-methylphenyl)propan-1-one (other names: 3-MMC; 3-methylmethcathinone)	1259
(105) 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)pentan-1-one (other names: dipentylone; *N*,*N*-dimethylpentylone)	7552
(106) 4-Chloromethcathinone (4-CMC, 1-(4-chlorophenyl)-2-(methylamino)propan-1-one)	1239
(107) 1-pentyl-3-(1-naphthoyl)indole (JWH-018 and AM678)	7118
(108) 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201)	7201
(109) 3-methoxyphencyclidine (Other names: 1-(1-(3-methoxyphenyl)cyclohexyl)piperidine; 3-MeO-PCP)	7457
(110) methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1*H*-indazole-3-carboxamido)butanoate (other name: MDMB-4en-PINACA)	7090
(111) Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate (other names: 4F-MDMB-BUTICA; 4F-MDMB-BICA)	7091
(112) *N*-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-en-1-yl)-1*H*-indazole-3-carboxamide (other name: ADB-4en-PINACA)	7092
(113) Ethyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate (other names: 5F-EDMB-PICA; 5F-EDMB-2201)	7094
(114) Methyl 2-(1-(4-fluorobenzyl)-1*H*-indole-3-carboxamido)-3-methyl butanoate (other name: MMB-FUBICA)	7095
(115) 6,6,9-trimethyl-3-pentyl-6a,7,8,9,10,10a-hexahydro-6*H*-benzo[*c*]chromen-1-ol (other names: hexahydrocannabinol, HHC)	7220
(116) 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido[4,3-b]indol-1-one (other names: CUMYL-PEGACLONE; SGT-151)	7093

(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) Clonazolam (Other name: 6-(2-chlorophenyl)-1-methyl-8-nitro-4*H*-benzo[*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepine)	2786
(2) Diclazepam (Other name: 7-chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2*H*-benzo[e][1,4]diazepin-2-one)	2789

(3) Etizolam (Other name: 4-(2-chlorophenyl)-2-ethyl-9-methyl-6*H*-thieno[3,2-*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepine)	2780
(4) Flualprazolam (Other name: 8-chloro-6-(2-fluorophenyl)-1-methyl-4*H*-benzo[*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepine)	2785
(5) Flubromazolam (Other name: 8-bromo-6-(2-fluorophenyl)-1-methyl-4*H*-benzo[*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepine)	2788
(6) gamma-hydroxybutyric acid (some other names include GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate)	2010
(7) Mecloqualone	2572
(8) Methaqualone	2565

(f) ***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:

(1) Amineptine (7-[(10,11-dihydro-5*H*-dibenzo[*a,d*]cyclohepten-5-yl)amino]heptanoic acid)	1219
(2) Aminorex (Some other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenly-2-oxazolamine)	1585
(3) N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine)	7493
(4) Cathinone	1235
Some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone, 2-aminopropiophenone, and norephedrone	
(5) 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
(6) Ethylphenidate (ethyl 2-phenyl-2-(piperidin-2-yl)acetate)	1727
(7) Fenethylline	1503
(8) 4-Fluoroamphetamine (4-FA, 1-(4-fluorophenyl)propan-2-amine, *para*-fluoroamphetamine)	1476
(9) Mesocarb (*N*-phenyl-*N* '-(3-(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-yl)carbamimidate)	1227
(10) Methcathinone (Some other names: 2-(methylamino)-propiophenone; alpha-(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-*N*-methylaminopropiophenone; monomethylpropion; ephedrone; *N*-methylcathinone; methylcathinone; AL-464; AL-422; AL-463 and UR1432), its salts, optical isomers and salts of optical isomers	1237
(11) Methiopropamine (*N*-methyl-1-(thiophen-2-yl)propan-2-amine)	1478
(12) (±)*cis*-4-methylaminorex ((±)*cis*-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine)	1590
(13) N-ethylamphetamine	1475
(14) *N,N*-dimethylamphetamine (also known as *N,N*-alpha-trimethyl-benzeneethanamine;	1480

N,N-alpha-trimethylphenethylamine)

(g) ***Cannabimimetic agents.*** Unless specifically exempted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of cannabimimetic agents, 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:

(1) Cannabimimetic agents 7000

(i) In this paragraph (g), *cannabimimetic agent* means any substance that is a cannabinoid receptor type 1 (CB1 receptor) agonist as demonstrated by binding studies and functional assays within any of the following structural classes:

(A) 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or alkenyl, whether or not substituted on the cyclohexyl ring to any extent.

(B) 3-(1-naphthoyl)indole or 3-(1-naphthylmethane)indole by substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the naphthoyl or naphthyl ring to any extent.

(C) 3-(1-naphthoyl)pyrrole by substitution at the nitrogen atom of the pyrrole ring, whether or not further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to any extent.

(D) 1-(1-naphthylmethylene)indene by substitution of the 3-position of the indene ring, whether or not further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to any extent.

(E) 3-phenylacetylindole or 3-benzoylindole by substitution at the nitrogen atom of the indole ring, whether or not further substituted in the indole ring to any extent, whether or not substituted on the phenyl ring to any extent.

(ii) The definition of cannabimimetic agent in this paragraph (g) includes, but is not limited to, the following substances:

(A) 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol (CP-47,497);

(B) 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol (cannabicyclohexanol or CP-47,497 C8-homolog);

(C) 1-butyl-3-(1-naphthoyl)indole (JWH-073);

(D) 1-hexyl-3-(1-naphthoyl)indole (JWH-019);

(E) 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWH-200);

(F) 1-pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250);

(G) 1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);

(H) 1-pentyl-3-(4-methyl-1-naphthoyl)indole (JWH-122);

(I) 1-pentyl-3-(4-chloro-1-naphthoyl)indole (JWH-398);

(J) 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694);

(K) 1-pentyl-3-[(4-methoxy)-benzoyl]indole (SR-19 and RCS-4);

(L) 1-cyclohexylethyl-3-(2-methoxyphenylacetyl)indole (SR-18 and RCS-8);

(M) 1-pentyl-3-(2-chlorophenylacetyl)indole (JWH-203);
(N) (1-((1-methylpiperidin-2-yl)methyl)-1*H*-indol-3-yl)(naphthalen-1-yl)methanone (AM-1220);
(O) (2-iodophenyl)(1-((1-methylpiperidin-2-yl)methyl)-1*H*-indol-3-yl)methanone (AM-2233);
(P) (4-ethylnaphthalen-1-yl)(1-(5-fluoropentyl)-1*H*-indol-3-yl)methanone (EAM-2201);
(Q) (4-methoxynaphthalen-1-yl)(2-methyl-1-pentyl-1*H*-indol-3-yl)methanone (JWH-098);
(R) 3-((4-methylnaphthalen-1-yl)methyl)-1-pentyl-1*H*-indole (JWH-184);
(S) (4-methylnaphthalen-1-yl)(1-(2-morpholinoethyl)-1*H*-indol-3-yl)methanone (JWH-193);
(T) (4-ethylnaphthalen-1-yl)(1-pentyl-1*H*-indol-3-yl)methanone (JWH-210);
(U) (1-(5-fluoropentyl)-1*H*-indol-3-yl)(4-methylnaphthalen-1-yl)methanone (MAM-2201);
(V) (2-methyl-1-pentyl-1*H*-indol-3-yl)(naphthalen-1-yl)methanone (JWH-007);
(W) naphthalen-1-yl(1-(pent-4-en-1-yl)-1*H*-indol-3-yl)methanone (JWH-022);
(X) (1-hexyl-5-phenyl-1*H*-pyrrol-3-yl)(naphthalen-1-yl)methanone (JWH-147);
(Y) 2-(3-methoxyphenyl)-1-(1-pentyl-1*H*-indol-3-yl)ethan-1-one (JWH-302);
(Z) (5-(2-fluorophenyl)-1-pentyl-1*H*-pyrrol-3-yl)(naphthalen-1-yl)methanone (JWH-307);
(AA) (4-fluoronaphthalen-1-yl)(1-pentyl-1*H*-indol-3-yl)methanone (JWH-412);
(BB) (5-methyl-3-(morpholinomethyl)-2,3-dihydro-[1,4]oxazino[2,3,4-*hi*]indol-6-yl)(naphthalen-1-yl)methanone (WIN 55,212-2);
(CC) 2-(5-hydroxy-2-(3-hydroxypropyl)cyclohexyl)-5-(2-methyloctan-2-yl)phenol (CP-55,940);
(DD) 2-(3-hydroxycyclohexyl)-5-(2-methylheptan-2-yl)phenol (CP-47,497 C6 homolog); and
(EE) 2-(3-hydroxycyclohexyl)-5-(2-methyldecan-2-yl)phenol (CP-47,497 C9 homolog).

(2) [Reserved]

(h) ***Temporary listing of substances subject to emergency scheduling.*** Any material, compound, mixture or preparation which contains any quantity of the following substances:

(1)-(29) [Reserved]

(30) 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 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, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;

(C) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, 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.

(ii) This definition includes, but is not limited to, the following substances:

(A)-(B) [Reserved]

(31)-(64) [Reserved]

(65)-(78) [Reserved]

(79) 2-(2-((2,3-dihydrobenzofuran-5-yl)methyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine, its isomers, esters, salts of isomers, esters and ethers (Other name: Ethyleneoxynitazene)

(80) 2-(2-(benzodioxol-5-ylmethyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine, its isomers, esters, ethers, salts isomers, esters and ethers (Other names: Methylenedioxynitazene; 3',4'-methylenedioxynitazene)

(81) 2-(2-(4-ethoxybenzyl)-5-methyl-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine, its isomers, esters, ethers, salts, and esters and ethers (Other name: 5-methyl etodesnitazene)

(82) 2-(2-(4-ethoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N*-ethylethan-1-amine, its isomers, esters, ethers, salts, and salts and ethers (Other name: *N*-desethyl etonitazene)

(83) *N*-ethyl-2-(5-nitro-2-(4-propoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine its isomers, esters, ethers, salts, and salts and ethers (Other name: *N*-desethyl protonitazene)

(84) 2-(2-(4-ethoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N,N*-dimethylethan-1-amine, its isomers, esters, ethers, salts, and esters and ethers (Other name: *N,N*-dimethylamino etonitazene)

(85) 2-(4-isopropoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole, its isomers, esters, ethers, salts, and salts esters and ethers (Other name: *N*-pyrrolidino isotonitazene)

(86) 8-bromo-1-methyl-6-phenyl-4*H*-benzo[*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepine, its salts, isomers, and salts of isomers (Oth bromazolam)

(87) 2-Fluorodeschloroketamine, its salts, isomers, and salts of isomers (other name: 2-(2-fluorophenyl)-2-(methylamino)c also known as 2-FDCK)

(88) *O*-Desmethyltramadol (other names: *O*-DSMT; desmetramadol; 3-[(1*R,*2*R*)-2-[(dimethylamino)methyl]-1-hydroxycycloh

(89) Methyl (*E*)-2-((1'*S,*6'*S,*7'S)-6'-ethyl-4-methoxy-3-oxo-3',5',6',7',8',8*a*'-hexahydro-2'*H*-spiro[indoline-2,1'-indolizine]-7'-yl)-3-m (commonly known as mitragynine pseudoindoxyl) including its isomers, esters, ethers, salts, and salts of isomers, esters, whenever the existence of such isomers, esters, ethers, and salts is possible. Since nomenclature of this substance is not standardized, compounds of this structure, regardless of numerical designation of atomic positions are covered

(90) Methyl (*E*)-2-((2*S,*3*S,*7*aS,*12*aR,*12*bS*)-3-ethyl-7*a*-hydroxy-8-methoxy-1,2,3,4,6,7,7*a,*12,12*a,*12*b*-decahydroindolo[2,3-*a*]quinolizin-2-yl)- (commonly known as MGM-15; also known as dihydro-7-hydroxymitragynine) including its isomers, esters, ethers, salts, a isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible. Since nomencla substance is not internationally standardized, compounds of this structure, regardless of numerical designation of atomic covered

(91) Methyl (*E*)-2-((2*S,*3*S,*7*aS,*12*aR,*12b*S*)-3-ethyl-9-fluoro-7*a*-hydroxy-8-methoxy-1,2,3,4,6,7,7*a,*12,12*a,*12*b*-decahydroindolo[2,3-*a*]quinolizin-2-yl)-3-methoxyacrylate (commonly known as MGM-16

9-fluoro-dihydro-7-hydroxymitragynine; or 10-fluoro-dihydro-7-hydroxymitragynine (depending on numbering convention)) i
isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, eth
possible. Since nomenclature of this substance is not internationally standardized, compounds of this structure, regardles
designation of atomic positions are covered

(92) 1-(1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-5,6-dichloro-1,3-dihydro-2*H*-benzo[*d*]imidazol-2-one, its isomers, esters, et
salts of isomers, esters, and ethers (Other names: 5,6-dichloro brorphine; SR-14968)

(93) 5,6-dichloro-1-(1-(4-chlorobenzyl)piperidin-4-yl)-1,3-dihydro-2*H*-benzo[*d*]imidazol-2-one, its isomers, esters, ethers, salt
isomers, esters, and ethers (Other names: 5,6-dichloro desmethylchlorphine; SR-17018)

(94) 3-(3-(1-(1-(4-chlorophenyl)ethyl)piperidin-4-yl)-2-oxo-2,3-dihydro-1*H*-benzo[*d*]imidazol-1-yl)propanenitrile, its isomers, e
salts, and salts of isomers, esters, and ethers (Other names: *N*-propionitrile chlorphine; cychlorphine)

(95) 8-(1-(4-chlorophenyl)ethyl)-1-phenyl-1,3,8-triazaspiro[4.5]decan-4-one, its isomers, esters, ethers, salts, and salts of is
ethers (Other names: spirochlorphine; R-6890)

[39 FR 22141, June 20, 1974]

Editorial Note: For FEDERAL REGISTER citations affecting § 1308.11, see the List of CFR Sections Affected, which appears in the Finding Aids section of the printed volume and at *www.govinfo.gov*.

Effective Date Notes:

1.At 83 FR 5191, Feb. 6, 2018, § 1308.11 was amended by adding paragraph (h)(30), effective Feb. 6, 2018, through Feb. 6, 2020. Effective Feb. 6, 2020, Congress extended the effective period for paragraph (h)(30) until May 6, 2021, by Public Law 116-114. Effective May 4, 2021, Congress extended the effective period for paragraph (h)(30) until October 22, 2021, by Public Law 117-12. Effective Sept. 30, 2021, Congress extended the effective period for paragraph (h)(30) until Jan. 28, 2022, by Public Law 117-43. Effective Jan. 13, 2022, Congress extended the effective period for paragraph (h)(30) until Feb. 18, 2022, by Public Law 117-70. Effective Feb. 18, 2022, Congress extended the effective period for paragraph (h)(30) until Mar. 11, 2022, by Public Law 117-86. Effective Mar. 11, 2022, Congress extended the effective period for paragraph (h)(30) until Mar. 15, 2022 by Public Law 117-95. Effective Mar. 15, 2022, Congress extended the effective period for paragraph (h)(30) until Dec. 31, 2022 by Public Law No. 117-103. Effective Dec. 29, 2022, Congress extended the effective period for paragraph (h)(30) until Dec. 31, 2024 by Public Law No. 117-328. Congress extended the effective period for paragraph (h)(30) until Mar. 31, 2025 by Public Law No. 116-114. Congress extended the effective period for (h)(30) until Sept. 30, 2025 by Public Law No. 119-4.

2. At 88 FR 86045, Dec. 12, 2023, § 1308.11 was amended by adding paragraphs (h)(62) through (h)(67), effective Dec. 12, 2023 through Dec. 12, 2025. At 90 FR 57356 and 57542, Dec. 11, 2025, paragraphs (h)(62) and (65) were extended through Dec. 12, 2026. At 91 FR 21963, Apr. 24, 2026, (h)(62) was removed.

3. At 90 FR 48265, Oct. 15, 2025, § 1308.11 was amended by adding paragraphs (h)(79) through (85), effective Oct. 15, 2025, until Oct. 15, 2027.

4. At 91 FR 12508, Mar. 16, 2026, § 1308.11 was amended by adding paragraph (h)(86), effective from Mar. 16, 2026, until Mar. 16, 2028.

5. At 91 FR 30209, May 22, 2026, § 1308.11 was amended by adding paragraph (h)(87), effective May 22, 2026 through May 22, 2028.

6. At 91 FR 54956, Aug. 26, 2026, § 1308.11 was amended by adding paragraphs (h)(88) through (h)(91), effective Aug. 26, 2026 through Aug. 26, 2028.

7. At 91 FR 55258, Aug. 27, 2026, § 1308.11 was amended by adding paragraphs (h)(92) through (h)(95), effective Aug. 27, 2026 through Aug. 27, 2028.