# 2023 Kansas Statutes

**65-4105. Substances included in schedule I.** (a) The controlled substances listed in this section are included in schedule I and the number set forth opposite each drug or substance is the DEA controlled substances code that has been assigned to it.
(b) Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers and salts is possible within the specific chemical designation:
(1) Acetyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N- phenylacetamide) 9821
(2) Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide) 9815
(3) Acetylmethadol 9601
(4) Acryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide; acryloylfentanyl) 9811
(5) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)cyclohexylmethyl]benzamide) 9551
(6) Allylprodine 9602
(7) Alphacetylmethadol 9603(except levo-alphacetylmethadol also known as levo-alpha-acetylmethadol, levomethadyl acetate or LAAM)
(8) Alphameprodine 9604
(9) Alphamethadol 9605
(10) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl] propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine) 9814
(11) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) 9832
(12) Benzethidine 9606
(13) Betacetylmethadol 9607
(14) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide) 9830
(15) Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide) 9831
(16) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide) 9836
(17) Betameprodine 9608
(18) Betamethadol 9609
(19) Betaprodine 9611
(20) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide) 9822
(21) Clonitazene 9612
(22) Crotonyl fentanyl ((E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide) 9844
(23) Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide) 9847
(24) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide) 9845
(25) Dextromoramide 9613
(26) Diampromide 9615
(27) Diethylthiambutene 9616
(28) Difenoxin 9168
(29) Dimenoxadol 9617
(30) Dimepheptanol 9618
(31) Dimethylthiambutene 9619
(32) Dioxaphetyl butyrate 9621
(33) Dipipanone 9622
(34) Ethylmethylthiambutene 9623
(35) Etonitazene 9624
(36) Etoxeridine 9625
(37) Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide) 9834
(38) Furethidine 9626

(39) Hydroxypethidine 9627
(40) Isotonitazene (N,N-diethyl-2-(2-(4 isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine; N,N-diethyl-2-[[4-(1-methylethoxy)phenyl]methyl]-5-nitro-1 H-benzimidazole-1-ethanamine) 9614
(41) Isobutyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide) 9827
(42) Ketobemidone 9628
(43) Levomoramide 9629
(44) Levophenacylmorphan 9631
(45) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide) 9825
(46) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide) 9813
(47) 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) 9833
(48) Morpheridine 9632
(49) Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide) 9838
(50) O-desmethyltramadolSome trade or other names: 2-((dimethylamino)methyl-1-(3-hydroxyphenyl)cyclohexanol;3-(2-((dimethylamino)methyl)-1-hydroxycyclohexyl)phenol
(51) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) 9661
(52) MT-45 (1-cychohexyl-4-(1,2-diphenylethyl)piperazine) 9560
(53) Noracymethadol 9633
(54) Norlevorphanol 9634
(55) Normethadone 9635
(56) Norpipanone 9636
(57) Ortho-fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide; 2-fluorofentanyl) 9816
(58) Para-chloroisobutyryl fentanyl (N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide) 9826
(59) Para-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) 9823
(60) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl]propanamide) 9812
(61) Para-fluoroisobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide, 4-fluoroisobutyryl fentanyl) 9824
(62) Para-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide) 9837
(63) PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine) 9663
(64) Phenadoxone 9637
(65) Phenampromide 9638
(66) Phenomorphan 9647
(67) Phenoperidine 9641
(68) Piritramide 9642
(69) Proheptazine 9643
(70) Properidine 9644
(71) Propiram 9649
(72) Racemoramide 9645
(73) Tetrahydrofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide) 9843
(74) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide) 9835
(75) Tilidine 9750
(76) Trimeperidine 9646
(77) U-47700 (3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide) 9547
(78) Valeryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide) 9840
(c) Any of the following opium derivatives, their salts, isomers and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible within the specific chemical designation:
(1) Acetorphine 9319

(2) Acetyldihydrocodeine 9051
(3) Benzylmorphine 9052
(4) Brorphine 9098
(5) Codeine methylbromide 9070
(6) Codeine-N-Oxide 9053
(7) Cyprenorphine 9054
(8) Desomorphine 9055
(9) Dihydromorphine 9145
(10) Drotebanol 9335
(11) Etorphine (except hydrochloride salt) 9056
(12) Heroin 9200
(13) Hydromorphinol 9301
(14) Methyldesorphine 9302
(15) Methyldihydromorphine 9304
(16) Morphine methylbromide 9305
(17) Morphine methylsulfonate 9306
(18) Morphine-N-Oxide 9307
(19) Myrophine 9308
(20) Nicocodeine 9309
(21) Nicomorphine 9312
(22) Normorphine 9313
(23) Pholcodine 9314
(24) Thebacon 9315
(d) Any material, compound, mixture or preparation that contains any quantity of the following hallucinogenic substances, their salts, isomers and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible within the specific chemical designation:
(1) Alpha-ethyltryptamine 7249Some 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 7391Some trade or other names: 4-bromo-2,5-dimethoxy-alpha-methylphenethylamine; 4-bromo-2,5-DMA.
(3) 2,5-dimethoxyamphetamine 7396Some trade or other names: 2,5-dimethoxy-alpha-methyl-phenethylamine; 2,5-DMA.
(4) 4-methoxyamphetamine 7411Some trade or other names: 4-methoxy-alpha-methylphene-thylamine; paramethoxyamphetamine; PMA.
(5) 5-methoxy-3,4-methylenedioxy-amphetamine 7401
(6) 4-methyl-2,5-dimethoxy-amphetamine 7395Some trade or other names: 4-methyl-2,5-dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP".
(7) 3,4-methylenedioxy amphetamine 7400
(8) 3,4-methylenedioxymethamphetamine (MDMA) 7405
(9) 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, and MDEA) 7404
(10) N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-alpha-methyl-3,4-(methylenedioxy) phenethylamine, and N-hydroxy MDA) 7402
(11) 3,4,5-trimethoxy amphetamine 7390
(12) Bufotenine 7433Some trade or other names: 3-(Beta-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine.
(13) Diethyltryptamine 7434Some trade or other names: N,N-Diethyltryptamine; DET.
(14) Dimethyltryptamine 7435Some trade or other names: DMT.
(15) Ibogaine 7260Some trade or other names: 7-Ethyl-6,6 Beta,7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido[1',2':1,2]azepino[5,4-b]indole; Tabernanthe iboga
(16) Lysergic acid diethylamide 7315
(17) Marijuana 7360
(18) Mescaline 7381
(19) Parahexyl 7374Some trade or other names: 3-Hexyl-l-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl.
(20) Peyote 7415Meaning 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.
(21) N-ethyl-3-piperidyl benzilate 7482
(22) N-methyl-3-piperidyl benzilate 7484
(23) Psilocybin 7437
(24) Psilocyn 7438Some trade or other names: Psilocin.
(25) Ethylamine analog of phencyclidine 7455Some trade or other names: N-ethyl-1-phenyl-cyclo-hexylamine; (1-phenylcyclohexyl)ethylamine; N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE.
(26) Pyrrolidine analog of phencyclidine 7458Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCPy; PHP.
(27) Thiophene analog of phencyclidine 7470Some trade or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 2-thienyl analog of phencyclidine; TPCP; TCP.
(28) 1-[1-(2-thienyl)-cyclohexyl] pyrrolidine 7473Some other names: TCPy.
(29) 2,5-dimethoxy-4-ethylamphetamine 7399Some trade or other names: DOET.
(30) Salvia divinorum or salvinorum A; all parts of the plant presently classified botanically as salvia divinorum, 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.
(31) Datura stramonium, commonly known as gypsum weed or jimson weed; all parts of the plant presently classified botanically as datura stramonium, 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.
(32) N-benzylpiperazine 7493Some trade or other names: BZP.
(33) 1-(3-[trifluoromethylphenyl])piperazineSome trade or other names: TFMPP.
(34) 4-Bromo-2,5-dimethoxyphenethylamine 7392
(35) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7), its optical isomers, salts and salts of optical isomers 7348
(36) Alpha-methyltryptamine (other name: AMT) 7432
(37) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT), its isomers,salts and salts of isomers 7439
(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-nitrophenyl)ethanamine (2C-N) 7521
(46) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P) 7524
(47) 5–methoxy–N,N–dimethyltryptamine (5–MeO–DMT) 7431Some trade or other names: 5–methoxy–3–[2–(dimethylamino) ethyl]indole.
(48) 2–(4–iodo–2,5–dimethoxyphenyl)–N–(2–methoxybenzyl)ethanamine 7538Some trade or other names: 25I–NBOMe; 2C–I–NBOMe; 25I; Cimbi–5.
(49) 2–(4–chloro–2,5–dimethoxyphenyl)–N–(2–methoxybenzyl)ethanamine 7537Some trade or other names: 25C–NBOMe; 2C–C–NBOMe; 25C; Cimbi–82.
(50) 2–(4–bromo–2,5–dimethoxyphenyl)–N–(2–methoxybenzyl)ethanamine 7536Some trade or other names: 25B–NBOMe; 2C–B–NBOMe; 25B; Cimbi–36.
(51) 2-(2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamineSome trade or other names: 25H-NBOMe.
(52) 2-(2,5-dimethoxy-4-methylphenyl)-N-(2-methoxybenzyl)ethanamineSome trade or other names: 25D-NBOMe; 2C-D-NBOMe.
(53) 2-(2,5-dimethoxy-4-nitrophenyl)-N-(2-methoxybenzyl)ethanamineSome trade or other names: 25N-NBOMe, 2C-N-NBOMe.
(54) 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1 H-pyrrolo[2,3-b]pyridine-3-carboxamide (5F-CUMYL-P7AICA) 7085
(e) Any material, compound, mixture or preparation that 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) EtizolamSome trade or other names: (4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f][1,2,4]triazolo[4,3-a][1,4]diazepine)
(2) Mecloqualone 2572
(3) Methaqualone 2565
(4) Gamma hydroxybutyric acid
(f) Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation that 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) Aminorex 1585Some other names: Aminoxaphen 2-amino-5-phenyl-2-oxazoline or 4,5-dihydro-5-phenyl-2-oxazolamine
(2) Fenethylline 1503
(3) N-ethylamphetamine 1475
(4) (+)cis-4-methylaminorex ((+)cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) 1590
(5) N,N-dimethylamphetamine (also known as N,N-alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine) 1480
(6) Cathinone (some other names: 2-amino-1-phenol-1-propanone, alpha-amino propiophenone, 2-amino propiophenone and norphedrone) 1235
(7) Substituted cathinonesAny compound, except bupropion or compounds listed under a different schedule, structurally derived from 2–aminopropan–1–one by substitution at the 1-position with either phenyl, naphthyl, or thiophene ring systems, whether or not the compound is further modified in any of the following ways:
(A) By substitution in the ring system to any extent with alkyl, alkylenedioxy, alkoxy, haloalkyl, hydroxyl, or halide substituents, whether or not further substituted in the ring system by one or more other univalent substituents;
(B) by substitution at the 3-position with an acyclic alkyl substituent;
(C) by substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or methoxybenzyl groups; or
(D) by inclusion of the 2-amino nitrogen atom in a cyclic structure.
(g) Any material, compound, mixture or preparation that contains any quantity of the following substances:
(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts and salts of isomers
(2) N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical isomers, salts and salts of isomers
(h) Any of the following cannabinoids, their salts, isomers and salts of isomers, unless specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible within the specific chemical designation:
(1) Tetrahydrocannabinols 7370Meaning tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity such as the following: Delta 1 cis or trans tetrahydrocannabinol, and their optical isomers Delta 6 cis or trans tetrahydrocannabinol, and their optical isomers Delta 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.), except tetrahydrocannabinols in any of the following:
(A) Industrial hemp, as defined in K.S.A. 2-3901, and amendments thereto;
(B) solid waste, as defined in K.S.A. 65-3402, and amendments thereto, and hazardous waste, as defined in K.S.A. 65-3430, and amendments thereto, if such waste is the result of the cultivation, production or processing of industrial hemp, as defined in K.S.A. 2-3901, and amendments thereto, and such waste contains a delta-9 tetrahydrocannabinol concentration of not more than 0.3%; or
(C) hemp products, as defined in K.S.A. 2-3901, and amendments thereto, unless otherwise deemed unlawful pursuant to K.S.A. 2-3908, and amendments thereto.

(2) NaphthylmethylindolesAny compound containing a 1H-indol-3-yl-(1-naphthyl)methane structure with substitution at the nitrogen atom of the indole group by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not further substituted on the indole group to any extent and whether or not substituted on the benzyl or naphthyl ring to any extent.
(3) NaphthoylpyrrolesAny compound containing a 3-(1-naphthoyl)pyrrole structure with substitution at the nitrogen atom of the pyrrole group by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not further substituted on the pyrrole group to any extent, whether or not substituted on the benzyl or naphthyl ring to any extent.
(4) NaphthylmethylindenesAny compound containing a naphthylmethylindene structure with substitution at the 3-position of the indene group by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not further substituted on the indene group to any extent, whether or not substituted on the benzyl or naphthyl ring to any extent.
(5) CyclohexylphenolsAny compound containing a 2-(3-hydroxycyclohexyl)phenol structure with substitution at the 5-position of the phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group whether or not substituted on the cyclohexyl ring to any extent.
(6) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-benzoxazin-6-yl]-1-napthalenylmethanone.Some trade or other names: WIN 55,212-2.
(7) 9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-olSome trade or other names: HU-210, HU-211.
(8) Indole-3-carboxylate estersAny compound containing a 1H-indole-3-carboxylate ester structure with the ester oxygen bearing a naphthyl, quinolinyl, isoquinolinyl or adamantyl group and substitution at the 1 position of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted on the indole ring to any extent and whether or not substituted on the naphthyl, quinolinyl, isoquinolinyl, adamantyl or benzyl groups to any extent.
(9) Indazole-3-carboxamidesAny compound containing a 1H-indazole-3-carboxamide structure with substitution at the nitrogen of the carboxamide by a naphthyl, quinolinyl, isoquinolinyl, adamantyl, benzyl, 1-amino-1-oxoalkan-2-yl or 1-alkoxy-1-oxoalkan-2-yl group and substitution at the 1 position of the indazole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted on the indazole ring to any extent and whether or not substituted on the naphthyl, quinolinyl, isoquinolinyl, adamantyl, 1-amino-1-oxoalkan-2-yl, 1-alkoxy-1-oxoalkan-2-yl or benzyl groups to any extent.
(10) Indole-3-carboxamidesAny compound containing a 1H-indole-3-carboxamide structure with substitution at the nitrogen of the carboxamide by a naphthyl, quinolinyl, isoquinolinyl, adamantyl, benzyl, 1-amino-1-oxoalkan-2-yl or 1-alkoxy-1-oxoalkan-2-yl group and substitution at the 1 position of the indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted on the indole ring to any extent and whether or not further substituted on the naphthyl, quinolinyl, isoquinolinyl, adamantyl, 1-amino-1-oxoalkan-2-yl, 1-alkoxy-1-oxoalkan-2-yl or benzyl groups to any extent.
(11) (1H-indazol-3-yl)methanonesAny compound containing a (1H-indazol-3-yl)methanone structure with the carbonyl carbon bearing a naphthyl group and substitution at the 1 position of the indazole ring by an alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl, or 2-(4-morpholinyl)ethyl group, whether or not further substituted on the indazole ring to any extent and whether or not substituted on the naphthyl or benzyl groups to any extent.
(12) (1H-indol-3-yl)methanonesAny compound containing a (1H-indol-3-yl)methanone

structure with the carbonyl carbon bearing a naphthyl, quinolinyl, isoquinolinyl, adamantyl, phenyl, benzyl or tetramethylcyclopropyl group and substitution at the 1 position of the indole ring by an alkyl,haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, benzyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or tetrahydropyranylmethyl group, whether or not further substituted on the indole ring to any extent and whether or not substituted on the naphthyl, quinolinyl, isoquinolinyl, adamantyl, phenyl, benzyl or tetramethylcyclopropyl groups to any extent.

**History:** L. 1972, ch. 234, § 5; L. 1982, ch. 269, § 2; L. 1985, ch. 220, § 1; L. 1986, ch. 241, § 1; L. 1987, ch. 244, § 1; L. 1989, ch. 200, § 1; L. 1991, ch. 199, § 1; L. 1992, ch. 174, § 1; L. 1993, ch. 70, § 1; L. 1994, ch. 54, § 1; L. 2001, ch. 171, § 3; L. 2008, ch. 124, § 1; L. 2010, ch. 7, § 1; L. 2011, ch. 83, § 3; L. 2013, ch. 67, § 1; L. 2014, ch. 79, § 1; L. 2015, ch. 27, § 1; L. 2016, ch. 95, § 2; L. 2017, ch. 57, § 4; L. 2018, ch. 101, § 1; L. 2019, ch. 37, § 13; L. 2022, ch. 99, § 3; June 9.