

SENATE BILL 2258

By Norris

AN ACT to amend Tennessee Code Annotated, Title 39, Chapter 17, Part 4; Title 41, Chapter 21, Part 2; Title 53 and Title 63, relative to controlled or addictive substances.

BE IT ENACTED BY THE GENERAL ASSEMBLY OF THE STATE OF TENNESSEE:

SECTION 1. Tennessee Code Annotated, Section 39-17-403(g), is amended by deleting the subsection in its entirety and substituting the following:

(g) The commissioner of mental health and substance abuse services, upon the agreement of the commissioner of health, in cooperation with the board of pharmacy, and in consultation with the director of the Tennessee bureau of investigation, shall revise and republish the schedules annually.

SECTION 2. Tennessee Code Annotated, Section 39-17-404, is amended by deleting the section in its entirety and substituting the following:

(a) The controlled substances listed or to be listed in the schedules in this part are included by whatever official, common, usual, chemical, or trade name designated.

(b) Notwithstanding any law to the contrary, the following are excluded from all schedules:

(1) Non-narcotic substances excluded under 21 CFR 1308.22, as amended;

(2) Chemical preparations exempted under 21 CFR 1308.24, as amended;

(3) Veterinary anabolic steroid implant products excluded under 21 CFR 1308.26, as amended;

(4) Prescription products exempted under 21 CFR 1308.32, as amended;

(5) Anabolic steroid products exempted under 21 CFR 1308.34, as amended; and

(6) Certain cannabis plant material, and products made from such material, that contain tetrahydrocannabinols and that are exempted under 21 CFR 1308.35, as amended.

SECTION 3. Tennessee Code Annotated, Section 39-17-406, is amended by deleting the section in its entirety and substituting the following:

(a) Schedule I consists of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section.

(b) Opiates, unless specifically excepted or unless listed in another schedule, means 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; provided, that for the purposes of subdivision (b)(48)(B)(xv), 3-Methylfentanyl, only, "isomer" includes the optical and geometric isomers:

- (1) Acetylmethadol;
- (2) Allylprodine;
- (3) Alphacetylmethadol (except levo-alphacetylmethadol, also known as levo-alpha-acetylmethadol; levomethadyl acetate; or LAAM);
- (4) Alphameprodine;
- (5) Alphamethadol;
- (6) Benzethidine;
- (7) Betacetylmethadol;
- (8) Betameprodine;
- (9) Betamethadol;
- (10) Betaprodine;

- (11) Clonitazene;
- (12) Dextromoramide;
- (13) Diampromide;
- (14) Diethylthiambutene;
- (15) Difenoxin;
- (16) Dimenoxadol;
- (17) Dimepheptanol;
- (18) Dimethylthiambutene;
- (19) Dioxaphetyl butyrate;
- (20) Dipipanone;
- (21) Ethylmethylthiambutene;
- (22) Etonitazene;
- (23) Etoxidine;
- (24) Furethidine;
- (25) Hydroxypethidine;
- (26) Ketobemidone;
- (27) Levomoramide;
- (28) Levophenacilmorphan;
- (29) Morpheridine;
- (30) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- (31) Noracymethadol;
- (32) Norlevorphanol;
- (33) Normethadone;
- (34) Norpipanone;
- (35) PEPAP (1-(2-phenylethyl)-4-phenyl-4-acetoxypiperidine);

- (36) Phenadoxone;
- (37) Phenampromide;
- (38) Phenomorphan;
- (39) Phenoperidine;
- (40) Piritramide;
- (41) Proheptazine;
- (42) Properidine;
- (43) Propiram;
- (44) Racemoramide;
- (45) Tilidine;
- (46) Trimeperidine;
- (47) U-47700; or
- (48) Fentanyl derivatives and analogues:

(A) Unless specifically excepted, listed in another schedule, or contained within a pharmaceutical product approved by the United States food and drug administration, any material, compound, mixture, or preparation, including its salts, isomers, esters, or ethers, and salts of isomers, esters, or ethers, whenever the existence of such salts is possible within any of the following specific chemical designations containing a 4-anilidopiperidine structure:

- (i) With or without substitution at the carbonyl of the aniline moiety with alkyl, alkenyl, carboalkoxy, cycloalkyl, methoxyalkyl, cyanoalkyl, or aryl groups, or furanyl, dihydrofuranyl, benzyl moiety, or rings containing heteroatoms sulfur, oxygen, or nitrogen;

(ii) With or without substitution at the piperidine amino moiety with a phenethyl, benzyl, alkylaryl (including heteroaromatics), alkyltetrazolyl ring, or an alkyl or carbomethoxy group, whether or not further substituted in the ring or group;

(iii) With or without substitution or addition to the piperidine ring to any extent with one or more methyl, carbomethoxy, methoxy, methoxymethyl, aryl, allyl, or ester groups;

(iv) With or without substitution of one or more hydrogen atoms for halogens, or methyl, alkyl, or methoxy groups, in the aromatic ring of the anilide moiety;

(v) With or without substitution at the alpha or beta position of the piperidine ring with alkyl, hydroxyl, or methoxy groups;

(vi) With or without substitution of the benzene ring of the anilide moiety for an aromatic heterocycle; or

(vii) With or without substitution of the piperidine ring for a pyrrolidine ring, perhydroazepine ring, or azepine ring; and

(B) The application of subdivision (b)(48)(A) includes, but is not limited to, any of the following:

(i) Acetylfentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);

(ii) Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidnyl]-N-phenyl-acetamide);

(iii) Acryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide);

(iv) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido)piperidine);

(v) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide);

(vi) Benzodioxolefentanyl;

(vii) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidiny]-N-phenylpropanamide);

(viii) Beta-hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide); N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidiny]-N-phenylpropanamide);

(ix) Beta-hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidiny]-N-phenylpropanamide);

(x) Butyrylfentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide; N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide);

(xi) Cyclopentyl fentanyl;

(xii) Isobutyryl fentanyl;

(xiii) Furanyl fentanyl;

(xiv) Lofentanil;

(xv) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide);

(xvi) 3-Methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide);

(xvii) Ocfentanil;

(xviii) Ohmefentanyl;

(xix) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide);

(xx) Para-fluoroisobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide; 4-fluoroisobutyryl fentanyl);

(xxi) Pentanoyl fentanyl;

(xxii) Thiofentanyl; or

(xxiii) Valeryl fentanyl.

(c) Opium derivatives, unless specifically excepted or unless listed in another schedule, means 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;

(2) Acetyldihydrocodeine;

(3) Benzylmorphine;

(4) Codeine methylbromide;

(5) Codeine-N-Oxide;

(6) Cyprenorphine;

(7) Desomorphine;

(8) Dihydromorphine;

(9) Drotebanol;

(10) Etorphine (except hydrochloride salt);

(11) Heroin;

- (12) Hydromorphenol;
- (13) Methyldesorphine;
- (14) Methyldihydromorphine;
- (15) Morphine methylbromide;
- (16) Morphine methylsulfonate;
- (17) Morphine-N-Oxide;
- (18) Myrophine;
- (19) Nicocodeine;
- (20) Nicomorphine;
- (21) Normorphine;
- (22) Pholcodine; or
- (23) Thebacon.

(d) Hallucinogenic substances, unless specifically excepted or unless listed in another schedule, means any material, compound mixture, or preparation that contains any quantity of the following hallucinogenic substances, or that 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 specified chemical designation; provided, that for purposes of this subsection (d) only, "isomer" includes the optical, positional, and geometric isomers:

(1) Alpha-ethyltryptamine

Other names: etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; [alpha]-ET; and AET; ET; Trip;

(2) Alpha-methyltryptamine

Other name: AMT;

(3) 4-Bromo-2,5-dimethoxyamphetamine

Other names: 4-Bromo-2,5-dimethoxy-[alpha]-methylphenethylamine; 4-bromo-2,5-DMA;

(4) 4-Bromo-2,5-dimethoxyphenethylamine

Other names: 2-(4-Bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl DOB; 2C-B; Nexus;

(5) 2-(4-Bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine

Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36;

(6) Bufotenine

Other names: 3-([beta]-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;

(7) 2-(4-Chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine

Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82;

(8) Diethyltryptamine

Other names: N,N-Diethyltryptamine; DET;

(9) 2,5-Dimethoxyamphetamine

Other names: 2,5-Dimethoxy-[alpha]-methylphenethylamine; 2,5-DMA;

(10) 2,5-Dimethoxy-4-ethylamphetamine

Other name: DOET;

(11) 2,5 Dimethoxy-4-(n)-propylthiophenethylamine

Other name: 2C-T-7;

(12) Dimethyltryptamine

Other name: DMT;

(13) Ethylamine analogue of phencyclidine

Other names: N-Ethyl-1-phenylcyclohexylamine; (1-phenylcyclohexyl)ethylamine;
N-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE;

(14) Ibogaine

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; Tabenanthe iboga;

(15) 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine

Other names: 25I-NBOMe; 2C-I-NBOMe; 25I; Cimbi-5;

(16) Lysergic acid diethylamide

Other name: LSD;

(17) Mescaline

Other name: Constituent of "Peyote" cacti;

(18) 4-Methoxyamphetamine

Other names: 4-methoxy-[alpha]-methylphenethylamine;
paramethoxyamphetamine; PMA;

(19) 5-Methoxy-3,4-methylenedioxyamphetamine;

(20) 5-Methoxy-N,N-diisopropyltryptamine

Other name: 5-MeO-DIPT;

(21) 5-methoxy-N,N-dimethyltryptamine

Other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT;

(22) 4-Methyl-2,5-dimethoxyamphetamine

Other names: 4-methyl-2,5-dimethoxy-[alpha]-methylphenethylamine; DOM;
STP;

(23) 3,4-Methylenedioxyamphetamine;

(24) 3,4-Methylenedioxymethamphetamine

Other name: MDMA;

(25) 3,4-Methylenedioxy-N-ethylamphetamine

Other names: N-ethyl-alpha-methyl 3,4(methylenedioxy) phenethylamine; N-ethyl MDA; MDE; MDEA;

(26) 3,4-Methylenedioxy-N-methylcathinone

Other name: Methylone;

(27) N-Ethyl-3-piperidyl benzilate;

(28) N-Hydroxy-3,4-methylenedioxyamphetamine

Other names: N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine; N-hydroxy MDA;

(29) N-methyl-3-piperidyl benzilate;

(30) Parahexyl

Other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl;

(31) Peyote

Meaning all parts of the plant presently classified botanically as *Lophophora williamsii* Lamaire, 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 or its seeds or extracts (Interprets 21 U.S.C. § 812(c), Schedule I(c)(12));

(32) Psilocybin (constituent of magic mushrooms);

(33) Psilocyn (constituent of magic mushrooms);

(34) Pyrrolidine analogue of phencyclidine (1-(1-phenylcyclohexyl)-pyrrolidine)

Other names: PCPy; PHP;

(35) 1-[1-(2-Thienyl)cyclohexyl]pyrrolidine

Other name: TCPy;

(36) 4-Methylmethcathinone

Other names: mephedrone; methpadrone; 4-MMC;

(37) 3,4-Methylenedioxyprovalerone

Other name: MDPV;

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

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

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

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

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

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

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

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

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

(47) Thiophene analogue of phencyclidine

Other names: 1-[1-(2-thienyl)cyclohexyl]piperidine; 2-thienylanalog of phencyclidine; TPCP; TCP;

(48) 3,4,5-Trimethoxyamphetamine;

(49) (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; or

(50) [1-(5-Fluoro-pentyl)-1H-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)

methanone

Other names: 5-fluoro-UR-144; 5-F-UR-144; XLR-11 1-(5-fluoro-pentyl)-3-(2,2,3,3-tetramethylcyclopropyl)indole.

(e) Depressants, unless specifically excepted or unless listed in another schedule, means 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 specified chemical designation:

(1) Etizolam

Other names: Etilaam, Etizola, Sedekopan, Pasaden, Depas;

(2) Gamma-hydroxybutyric acid

Other names: GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate;

(3) Mecloqualone; or

(4) Methaqualone.

(f) Stimulants, unless specifically excepted or unless listed in another schedule, means 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) Alpha-pyrrolidinobutiophenone

Other names: [alpha]-PBP; 1-phenyl-2-(pyrrolidin-1-yl)butan-1-one;

(2) Alpha-pyrrolidinopentiophenone

Other names: [alpha]-PVP; [alpha]-pyrrolidinovalerophenone; 1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one;

(3) Aminorex

Other names: aminoxophen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine;

(4) Butylone

Other names: bk-MBDB; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one;

(5) Cathinone

Other names: 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone; 2-aminopropiophenone; norphedrone; constituent of *catha edulis* or "Khat" plant;

(6) 3-Fluoro-N-methylcathinone

Other names: 3-FMC; 1-(3-fluorophenyl)-2-(methylamino)propan-1-one;

(7) 4-Fluoro-N-methylcathinone

Other names: 4-FMC; flephedrone; 1-(4-fluorophenyl)-2-(methylamino)propan-1-one;

(8) Fenethylamine;

(9) Methcathinone

Other names: 2-(methylamino)-propionophenone; alpha-(methylamino)propionophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-N-methylaminopropionophenone; monomethylpropion; ephedrone; N-methylcathinone; methylcathinone; AL-464; AL-422; AL-463; and UR1432;

(10) (+/-)-cis-4-Methylaminorex (cis isomer)

Other name: (+/-)-cis-4,5 dihydro-4-methyl-5-phenyl-2-oxazolamine;

(11) 4-Methyl-N-ethylcathinone

Other names: 4-MEC; 2-(ethylamino)-1-(4-methylphenyl)propan-1-one;

(12) 4-Methyl-alpha-pyrrolidinopropiophenone

Other names: 4-MePPP; MePPP; 4-methyl-[alpha]-pyrrolidinopropiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)-propan-1-one;

(13) Naphyrone

Other names: naphthylpyrovalerone; 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one;

(14) N-Benzylpiperazine

Other names: BZP; 1-benzylpiperazine;

(15) N-Ethylamphetamine;

(16) N,N-Dimethylamphetamine

Other names: N,N-alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine;

(17) Pentedrone

Other names: [alpha]-methylaminovalerophenone; 2-(methylamino)-1-phenylpentan-1-one; or

(18) Pentylone

Other names: bk-MBDP; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one.

(g) Cannabimimetic agents, unless specifically exempted or unless listed in another schedule, means any material, compound, mixture, or preparation that contains any quantity of the following substances, or that 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) 5-(1,1-Dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol (CP-47,497);

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

(3) 1-Pentyl-3-(1-naphthoyl)indole (JWH-018 and AM678);

(4) 1-Butyl-3-(1-naphthoyl)indole (JWH-073);

(5) 1-Hexyl-3-(1-naphthoyl)indole (JWH-019);

(6) 1-[2-(4-Morpholinyl)ethyl]-3-(1-naphthoyl)indole (JWH-200);

(7) 1-Pentyl-3-(2-methoxyphenylacetyl)indole (JWH-250);

(8) 1-Pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);
(9) 1-Pentyl-3-(4-methyl-1-naphthoyl)indole (JWH-122);
(10) 1-Pentyl-3-(4-chloro-1-naphthoyl)indole (JWH-398);
(11) (1-(5-Fluoropentyl)-1H-indazol-3-yl)(naphthalen-1-yl)methanone
(THJ-2201);

(12) 1-(5-Fluoropentyl)-3-(1-naphthoyl)indole (AM2201);
(13) 1-(5-Fluoropentyl)-3-(2-iodobenzoyl)indole (AM694);
(14) 1-Pentyl-3-[(4-methoxy)-benzoyl]indole (SR-19 and RCS-4);
(15) 1-Cyclohexylethyl-3-(2-methoxyphenylacetyl)indole (SR-18 and

RCS-8);

(16) 1-Pentyl-3-(2-chlorophenylacetyl)indole (JWH-203);
(17) Methyl 2-(1-(cyclohexylmethyl)-1h-indole-3-carboxamido)-3,3-

dimethylbutanoate

Other names: MDMB-CHMICA, MMB-CHMINACA;

(18) Methyl 2-(1-(4-fluorobenzyl)-1h-indazole-3-carboxamido)-3,3-
dimethylbutanoate

Other name: MDMB-FUBINACA;

(19) Methyl 2-(1-(5-fluoropentyl)-1h-indazole-3-carboxamido)-3-
methylbutanoate

Other name: 5F-AMB;

(20) Methyl 2-(1-(5-fluoropentyl)-1h-indazole-3-carboxamido)-3,3-
dimethylbutanoate

Other names: 5F-ADB, 5F-MDMB-PINACA;

(21) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide

Other names: APINACA; AKB48;

(22) N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide

Other names: 5F-APINACA, 5F-AKB48;

(23) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide

Other name: AB-FUBINACA;

(24) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide

Other name: AB-CHMINACA;

(25) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)1H-indazole-3-carboxamide

Other name: ADB-FUBINACA;

(26) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide

Other name: ADB-PINACA;

(27) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide

Other name: AB-PINACA;

(28) Quinolin-8-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate

Other names: 5-fluoro-PB-22; 5F-PB-22; or

(29) Quinolin-8-yl 1-pentyl-1H-indole-3-carboxylate

Other names: PB-22; QUPIC.

SECTION 4. Tennessee Code Annotated, Section 39-17-408(c), is amended by deleting the subsection in its entirety and substituting the following:

(c) Opiates, unless specifically excepted or unless in another schedule, means 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:

- (1) Alfentanil;
- (2) Alphaprodine;
- (3) Anileridine;
- (4) Bezitramide;
- (5) Carfentanil;
- (6) Dextropropoxyphene (bulk, non-dosage forms);
- (7) Dihydrocodeine;
- (8) Diphenoxylate;
- (9) Fentanyl;
- (10) Isomethadone;
- (11) Levo-alpha-acetylmethadol

Other names: levo-alpha-acetylmethadol; levomethadyl acetate; LAAM;

- (12) Levomethorphan;
- (13) Levorphanol;
- (14) Metazocine;
- (15) Methadone;
- (16) Methadone-Intermediate; 4-cyano-2-dimethylamino-4,4-diphenyl butane;
- (17) Moramide-Intermediate; diphenylpropane-carboxylic acid; 2-methyl-3-morpholino-1,1-Pethidine (meperidine);
- (18) Pethidine (meperidine);
- (19) Pethidine-Intermediate-A; 4-cyano-1-methyl-4-phenylpiperidine;

- (20) Pethidine-Intermediate-B; ethyl-4-phenylpiperidine-4-carboxylate;
- (21) Pethidine-Intermediate-C; 1-methyl-4-phenylpiperidine-4-carboxylic acid;
- (22) Phenazocine;
- (23) Piminodine;
- (24) Racemethorphan;
- (25) Racemorphan;
- (26) Remifentanil;
- (27) Sufentanil;
- (28) Tapentadol; or
- (29) Thiafentanil.

SECTION 5. Tennessee Code Annotated, Section 39-17-408(f), is amended by deleting the subsection in its entirety and substituting the following:

(f) Hallucinogenic substances:

(1) Nabilone

Other names: (+/-)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1-hydroxy-6,6-dimethyl-9H-dibenzo[b,d]pyran-9-one; or

(2) Dronabinol in oral solution in drug product approved for marketing by United States food and drug administration

Other names: [(-)-delta-9-trans tetrahydrocannabinol], Syndros.

SECTION 6. Tennessee Code Annotated, Section 39-17-408(g)(2)(A), is amended by deleting the subdivision in its entirety and substituting the following:

(A) 1-phenylcyclohexylamine; or

SECTION 7. Tennessee Code Annotated, Section 39-17-410(f)(1), is amended by deleting the subdivision in its entirety and substituting the following:

(1) Anabolic steroids:

- (A) 3[alpha],17[beta]-dihydroxy-5a-androstane;
- (B) 17[alpha]-methyl-3[alpha],17[beta]-dihydroxy-5a-androstane;
- (C) 17[alpha]-methyl-3[beta],17[beta]-dihydroxy-5a-androstane;
- (D) 17[alpha]-methyl-3[beta],17[beta]-dihydroxyandrost-4-ene;
- (E) 17[alpha]-methyl-[delta]1-dihydrotestosterone(17[beta]-hydroxy-

17[alpha]-methyl-5[alpha]-androst-1-en-3-one

Other Names: 17-[alpha]-methyl-1-testosterone;

(F) 17[alpha]-methyl-4-hydroxynandrolone(17[alpha]-methyl-4-hydroxy-17[beta]-hydroxyestr-4-en-3-one);

(G) 1-Androstenediol (3[alpha],17[beta]-dihydroxy-5[alpha]-androst-1-ene);

(H) 1-Androstenediol (3[beta],17[beta]-dihydroxy-5[alpha]-androst-1-ene);

(I) 4-Androstenediol (3[beta],17[beta]-dihydroxy-androst-4-ene);

(J) 5-Androstenediol (3[beta],17[beta]-dihydroxy-androst-5-ene);

(K) 1-Androstenedione (5[alpha]-androst-1-en-3,17-dione);

(L) 4-Androstenedione (androst-4-en-3,17-dione);

(M) 5-Androstenedione (androst-5-en-3,17-dione);

(N) 3[Beta],17[beta]-dihydroxy-5a-androstane;

(O) 13[Beta]-ethyl-17[beta]-hydroxygon-4-en-3-one;

(P) Androstenedione (5[alpha]-androst-1-en-3,17-dione);

(Q) Bolasterone (7[alpha],17[alpha]-dimethyl-17[beta]-hydroxyandrost-4-en-3-one);

(R) Boldenone (17[beta]-hydroxyandrost-1,4-diene-3-one);

(S) Boldione (androsta-1,4-diene-3,17-dione);

(T) Calusterone (7[beta],17[alpha]-dimethyl-17[beta]-hydroxyandrost-4-en-3-one);

(U) Clostebol (4-chloro-17[beta]-hydroxyandrost-4-en-3-one)

Other Name: 4-Chlorotestosterone;

(V) Dehydrochloromethyltestosterone (4-chloro-17[beta]-hydroxy-17[alpha]-methylandrost-1,4-dien-3-one);

(W) [Delta]1-dihydrotestosterone (17[beta]-hydroxy-5[alpha]-androst-1-en-3-one)

Other name: 1-testosterone;

(X) Desoxymethyltestosterone (17[alpha]-methyl-5[alpha]-androst-2-en-17[beta]-ol)

Other name: madol;

(Y) 4-Dihydrotestosterone (17[beta]-hydroxyandrostan-3-one);

(Z) Drostanolone (17[beta]-hydroxy-2[alpha]-methyl-5[alpha]-androstan-3-one);

(AA) Ethylestrenol (17[alpha]-ethyl-17[beta]-hydroxyestr-4-ene);

(BB) Fluoxymesterone (9-fluoro-17[alpha]-methyl-11[beta],17[beta]-dihydroxyandrost-4-en-3-one);

(CC) Formebolone (2-formyl-17[alpha]-methyl-11[alpha],17[beta]-dihydroxyandrost-1,4-dien-3-one);

(DD) Furazabol (17[alpha]-methyl-17[beta]-hydroxyandrostano[2,3-c]-furazan);

(EE) 4-Hydroxy-19-nortestosterone (4,17[beta]-dihydroxyestr-4-en-3-one);

(FF) 4-Hydroxytestosterone (4,17[beta]-dihydroxyandrost-4-en-3-one);

(GG) Mestanolone (17[alpha]-methyl-17[beta]-hydroxy-5[alpha]-androstan-3-one);

(HH) Mesterolone (1[alpha]-methyl-17[beta]-hydroxy-5[alpha]-androstan-3-one);

(II) Methandienone (17[alpha]-methyl-17[beta]-hydroxyandrost-1,4-diene-3-one);

(JJ) Methandranone;

(KK) Methandriol (17[alpha]-methyl-3[beta],17[beta]-dihydroxyandrost-5-ene);

(LL) Methandrostenolone;

(MM) Methasterone (2[alpha],17[alpha]-dimethyl-5[alpha]-androstan-17[beta]-ol-3-one);

(NN) Methenolone (1-methyl-17[beta]-hydroxy-5[alpha]-androst-1-en-3-one);

(OO) Methyldienolone (17[alpha]-methyl-17[beta]-hydroxyestra-4,9(10)-dien-3-one);

(PP) Methyltestosterone (17[alpha]-methyl-17[beta]-hydroxyandrost-4-en-3-one);

(QQ) Methyltrienolone (17[alpha]-methyl-17[beta]-hydroxyestra-4,9,11-trien-3-one);

(RR) Mibolerone (7[alpha],17[alpha]-dimethyl-17[beta]-hydroxyestr-4-en-3-one);

(SS) Nandrolone (17[beta]-hydroxyestr-4-en-3-one);

(TT) 19-Nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione);

(UU) 19-Nor-4-androstenediol (3[alpha],17[beta]-dihydroxyestr-4-ene);

(VV) 19-Nor-4-androstenediol (3[beta],17[beta]-dihydroxyestr-4-ene);

(WW) 19-Nor-5-androstenediol (3[alpha],17[beta]-dihydroxyestr-5-ene);

(XX) 19-Nor-5-androstenediol (3[beta],17[beta]-dihydroxyestr-5-ene);

(YY) 19-Nor-4-androstenedione (estr-4-en-3,17-dione);

(ZZ) 19-Nor-5-androstenedione (estr-5-en-3,17-dione);

(AAA) Norbolethone (13[beta],17[alpha]-diethyl-17[beta]-hydroxygon-4-en-3-one);

(BBB) Norclostebol (4-chloro-17[beta]-hydroxyestr-4-en-3-one);

(CCC) Norethandrolone (17[alpha]-ethyl-17[beta]-hydroxyestr-4-en-3-one);

(DDD) Normethandrolone (17[alpha]-methyl-17[beta]-hydroxyestr-4-en-3-one);

(EEE) Oxandrolone (17[alpha]-methyl-17[beta]-hydroxy-2-oxa-5[alpha]-androstan-3-one);

(FFF) Oxymesterone (17[alpha]-methyl-4,17[beta]-dihydroxyandrost-4-en-3-one);

(GGG) Oxymetholone (17[alpha]-methyl-2-hydroxymethylene-17[beta]-hydroxy-[5[alpha]]-androstan-3-one);

(HHH) Prostanazol (17[beta]-hydroxy-5[alpha]-androstan[3,2-c]pyrazole);

(III) Stanolone (17[beta]-hydroxy-5alpha-androstan-3-one);

(JJJ) Stanozolol (17[alpha]-methyl-17[beta]-hydroxy-[5[alpha]]-androst-2-eno[3,2-c]-pyrazole);

(KKK) Stenbolone (17[beta]-hydroxy-2-methyl-[5[alpha]]-androst-1-en-3-one);

(LLL) Testolactone (13-hydroxy-3-oxo-13,17-secoandrosta-1,4-dien-17-oic acid lactone);

(MMM) Testosterone (17[beta]-hydroxyandrost-4-en-3-one);

(NNN) Tetrahydrogestrinone (13[beta],17[alpha]-diethyl-17[beta]-hydroxygon-4,9,11-trien-3-one); or

(OOO) Trenbolone (17[beta]-hydroxyestr-4,9,11-trien-3-one).

SECTION 8. Tennessee Code Annotated, Section 39-17-412(g), is amended by deleting the subsection in its entirety and substituting the following:

(g) Other substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation that contains any quantity of the following substances, including its salts:

(1) Pentazocine;

(2) Butorphanol (including its optical isomers); or

(3) Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-oxopropyl]][(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and salts of isomers.

SECTION 9. Tennessee Code Annotated, Section 39-17-414(d), is amended by deleting the subsection in its entirety and substituting the following:

(d) Depressants, unless specifically exempted or excluded or unless listed in another schedule, means 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:

- (1) Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide);
 - (2) Ezogabine [N-[2-amino-4-(4-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester];
 - (3) Gabapentin [1-(aminomethyl)cyclohexaneacetic acid];
 - (4) Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide];
- or
- (5) Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid].

SECTION 10. Tennessee Code Annotated, Section 39-17-452(a)(2)(H), is amended by deleting the subdivision in its entirety.

SECTION 11. Tennessee Code Annotated, Section 39-17-452(a), is amended by adding the following language as a new subdivision (a)(1) and redesignating the existing subdivisions accordingly:

- (1) Unless specifically excepted or unless listed in another schedule, it is an offense to knowingly produce, manufacture, distribute, sell, offer for sale, or possess any capsule, pill, or other product composed of or containing any amount of Kratom, in its natural botanical form, or any capsule, pill, or other product composed of or containing any amount of mitragynine or hydroxymitragynine.

SECTION 12. Tennessee Code Annotated, Section 41-21-236(a)(2), is amended by redesignating subdivision (a)(2)(D) as subdivision (a)(2)(B)(iv), adding the following as a new subdivision (a)(2)(C), and redesignating the existing subdivisions accordingly:

(C)

(i) In addition to the credits authorized by subdivisions (a)(2)(A) and (B), the department shall provide a credit of sixty (60) days to any qualifying prisoner who successfully completes an evidence-based, intensive residential substance use disorder treatment therapeutic community program of at least nine (9) months in the department.

(ii) It is the legislative intent that this credit be implemented by the department in a manner that maximizes the potential of prisoners who have a history of substance use disorders returning to the community to become working and productive members of society by breaking the cycle of substance use and criminal behavior through successful completion of intensive substance use disorder treatment.

(iii) The provisions of this subdivision (a)(2)(C) permitting an additional sixty (60) days of credit for successful completion of the intensive substance use disorder treatment program shall not apply to any prisoner convicted of an offense that requires service of at least eighty-five percent (85%) of the sentence under § 40-35-501(i) or one hundred percent (100%) of the sentence under § 39-13-523 or § 40-35-501(l).

(iv) The sixty (60) days of credit authorized by this subdivision (a)(2)(C) shall apply to any inmate who successfully completes an evidence-based, intensive residential substance use disorder treatment therapeutic community program of at least nine (9) months in the department on or after July 1, 2018.

SECTION 13. If any provision of this act or the application of any provision of this act to any person or circumstance is held invalid, the invalidity does not affect other provisions or applications of the act that can be given effect without the invalid provision or application, and to that end, the provisions of this act are declared to be severable.

SECTION 14. This act shall take effect July 1, 2018, the public welfare requiring it.