

SCHEDULES OF CONTROLLED SUBSTANCES

PURSUANT TO THE TEXAS CONTROLLED SUBSTANCES ACT, HEALTH AND SAFETY CODE, CHAPTER 481, THESE SCHEDULES SUPERCEDE PREVIOUS SCHEDULES AND CONTAIN THE MOST CURRENT VERSION OF THE SCHEDULES OF ALL CONTROLLED SUBSTANCES FROM THE PREVIOUS SCHEDULES AND MODIFICATIONS.

This annual publication of the Texas Schedules of Controlled Substances was signed by John Hellerstedt, M.D., Commissioner of Health, and will take effect 21 days following publication of this notice in the *Texas Register*.

Changes to the schedules are designated by an asterisk (*). Additional information can be obtained by contacting the Department of State Health Services, Drugs and Medical Devices Unit, P.O. Box 149347, Austin, Texas 78714-9347. The telephone number is (512) 834-6755 and the website address is <http://www.dshs.texas.gov/dmd>.

SCHEDULES

Nomenclature: Controlled substances listed in these schedules are included by whatever official, common, usual, chemical, or trade name they may be designated.

SCHEDULE I

Schedule I consists of:

-Schedule I opiates

The following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, if the existence of these isomers, esters, ethers, and salts are possible within the specific chemical designation:

- (1) Acetyl- α -methylfentanyl (*N*-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-*N*-phenylacetamide);
- (2) Acetylmethadol;
- (3) Acetyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacetamide);
- (4) Acryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacrylamide) (Other name: acryloylfentanyl);
- (5) AH-7921 (3,4-dichloro-*N*-[1-(dimethylamino)cyclohexymethyl]benzamide);
- (6) Allylprodine;
- (7) Alphacetylmethadol (except levo- α -cetylmethadol, levo- α -acetylmethadol, levomethadyl acetate, or LAAM);
- (8) α -Methylfentanyl or any other derivative of fentanyl;

- (9) α -Methylthiofentanyl (*N*-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl] *N*-phenylpropanamide);
- (10) Benzethidine;
- (11) β -Hydroxyfentanyl (*N*-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-*N*-phenylpropanamide);
- (12) β -Hydroxy-3-methylfentanyl (*N*-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-*N*-phenylpropanamide);
- (13) β -hydroxythiofentanyl (Other names: *N*-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-*N*-phenylproprionamide; *N*-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidnyl]-*N*-phenylpropanamide);
- (14) Betaprodine;
- (15) Butyryl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylbutanamide);
- (16) Clonitazene;
- * (17) Crotonyl fentanyl (Other name: (6-2-5) (E)-*N*-(1-Phenethylpiperidin-4-yl)-*N*-phenylbut-2-enamide);
- (18) Cyclopropyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylcyclopropanecarboxamide);
- (19) Diampromide;
- (20) Diethylthiambutene;
- (21) Difenoxin;
- (22) Dimenoxadol;
- (23) Dimethylthiambutene;
- (24) Dioxaphetyl butyrate;
- (25) Dipipanone;
- (26) Ethylmethylthiambutene;
- (27) Etonitazene;
- (28) Etoxeridine;
- (29) 4-Fluoroisobutyryl fentanyl (*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide) (Other name: *p*-fluoroisobutyryl fentanyl);
- (30) Furanyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylfuran-2-carboxamide);
- (31) Furethidine;
- (32) Hydroxypethidine;
- (33) Ketobemidone;
- (34) Levophenacymorphan;
- (35) Meprodine;
- (36) Methadol;
- (37) Methoxyacetyl fentanyl (2-methoxy-*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylacetamide);
- (38) 3-Methylfentanyl (*N*-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-*N*-phenylpropanamide);
- (39) 3-Methylthiofentanyl (*N*-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-*N*-phenylpropanamide);
- (40) Moramide;
- (41) Morpheridine;
- (42) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- (43) MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
- (44) Noracymethadol;
- (45) Norlevorphanol;

- (46) Normethadone;
- (47) Norpipanone;
- (48) Ocfentanil (*N*-(2-fluorophenyl)-2-methoxy-*N*-(1-phenethylpiperidin-4-yl)acetamide);
- (49) *o*-Fluorofentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)propionamide) (Other name: 2-fluorofentanyl);
- (50) *p*-Fluorobutyryl fentanyl (*N*-(4-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)butyramide);
- (51) *p*-Fluorofentanyl (*N*-(4-fluorophenyl)-*N*-[1-(2-phenethyl)-4 piperidinyl]propanamide);
- (52) PEPAP (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- (53) Phenadoxone;
- (54) Phenampromide;
- (55) Phencyclidine;
- (56) Phenomorphan;
- (57) Phenoperidine;
- (58) Piritramide;
- (59) Proheptazine;
- (60) Properidine;
- (61) Propiram;
- (62) Tetrahydrofuranfentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenyltetrahydrofuran-2-carboxamide);
- (63) Thiofentanyl (*N*-phenyl-*N*-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);
- (64) Tilidine;
- (65) Trimeperidine; and
- (66) U-47700 (3,4-dichloro-*N*-[2-(dimethylamino)cyclohexyl]-*N*-methylbenzamide).

-Schedule I opium derivatives

The following opium derivatives, their salts, isomers, and salts of isomers, unless specifically excepted, if the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

The following opium derivatives, their salts, isomers, and salts of isomers, unless specifically excepted, if the existence of these 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) Hydromorphinol;

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

-Schedule I hallucinogenic substances

Unless specifically excepted or unless listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following hallucinogenic substances or that contains any of the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation (for the purposes of this Schedule I hallucinogenic substances section only, the term "isomer" includes optical, position, and geometric isomers):

- (1) α -Ethyltryptamine (Other names: etryptamine; Monase; α -ethyl-1*H*-indole-3-ethanamine; 3-(2-aminobutyl) indole; α -ET; AET);
- (2) 4-Bromo-2,5-dimethoxyamphetamine (Other names: 4-bromo-2,5- dimethoxy- α -methylphenethylamine; 4-bromo-2,5-DMA);
- (3) 4-Bromo-2,5-dimethoxyphenethylamine (Other names: Nexus; 2C-B; 2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; α -desmethyl DOB);
- (4) 2,5-Dimethoxyamphetamine (Other names: 2,5-dimethoxy- α -methylphenethylamine; 2,5-DMA);
- (5) 2,5-Dimethoxy-4-ethylamphetamine (Other name: DOET);
- (6) 2,5-Dimethoxy-4-(*n*)-propylthiophenethylamine, its optical isomers, salts and salts of isomers (Other name: 2C-T-7);
- (7) 4-Methoxyamphetamine (Other names: 4-methoxy- α -methylphenethylamine; paramethoxyamphetamine; PMA);
- (8) 5-Methoxy-3,4-methylenedioxyamphetamine (Other names: MMDA);
- (9) 4-Methyl-2,5-dimethoxyamphetamine (Other names: 4-methyl-2,5-dimethoxy- α -methyl-phenethylamine; "DOM"; "STP");
- (10) 3,4-Methylenedioxyamphetamine (Other names: MDA; Love Drug);
- (11) 3,4-Methylenedioxymethamphetamine (Other names: MDMA; MDM; Ecstasy; XTC);
- (12) 3,4-Methylenedioxy-*N*-ethylamphetamine (Other names: *N*-ethyl- α -methyl-3,4(methylenedioxy)phenethylamine; *N*-ethyl MDA; MDE; MDEA);
- (13) *N*-Hydroxy-3,4-methylenedioxyamphetamine (Other names: *N*-hydroxy MDA);
- (14) 3,4,5-Trimethoxyamphetamine (Other names: TMA);

(15) 5-Methoxy-*N,N*-dimethyltryptamine (Other names: 5-methoxy-3-[2-(dimethylamino)ethyl]indole, 5-MeO-DMT);

(16) α -Methyltryptamine (Other name: AMT), its isomers, salts, and salts of isomers;

(17) Bufotenine (Other names: 3- β -Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; *N,N*-dimethylserotonin; 5-hydroxy-*N,N*-dimethyltryptamine; mappine);

(18) Diethyltryptamine (Other names: *N,N*-Diethyltryptamine; DET);

(19) Dimethyltryptamine (Other name: DMT);

(20) 5-Methoxy-*N,N*-diisopropyltryptamine, its isomers, salts, and salts of isomers (Other name: 5-MeO-DIPT);

(21) Ibogaine (Other names: 7-Ethyl-6,6- β -7,8,9,10,12,13-octhydro-2-methoxy-6,9-methano-5*H*-pyrido[1',2':1,2] azepino [5,4-*b*] indole; Tabernanthe iboga);

(22) Lysergic acid diethylamide;

(23) Marihuana.

The term marihuana does not include hemp, as defined in Title 5, Agriculture Code, Chapter 121.

(24) Mescaline;

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

(26) Peyote, unless unharvested and growing in its natural state, meaning all parts of the plant classified botanically as *Lophophora williamsii* *Lemaire*, whether growing or not, the seeds of the plant, an extract from a part of the plant, and every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or extracts;

(27) *N*-ethyl-3-piperidyl benzilate;

(28) *N*-methyl-3-piperidyl benzilate;

(29) Psilocybin;

(30) Psilocyn;

*(31) Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained in a plant of the genus *Cannabis* (*cannabis* plant), except for up to 0.3 percent delta-9-tetrahydrocannabinols in hemp (as defined under Texas Agriculture Code 121), 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.)

(32) Ethylamine analog of phencyclidine (Other names: *N*-ethyl-1-phenylcyclohexylamine; (1-phenylcyclohexyl)ethylamine; *N*-(1-phenylcyclohexyl)ethylamine; cyclohexamine; PCE);

- (33) Pyrrolidine analog of phencyclidine (Other names: 1-(1 phenyl-cyclohexyl)-pyrrolidine; PCPy; PHP; rolicyclidine);
- (34) Thiophene analog of phencyclidine (Other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 2-thienyl analog of phencyclidine; TPCP; TCP);
- (35) 1-[1-(2-Thienyl)cyclohexyl]pyrrolidine (Other name: TCPy);
- (36) 4-Methylmethcathinone (Other names: 4-methyl-*N*-methylcathinone; mephedrone);
- (37) 3,4-methylenedioxypropylvalerone (MDPV);
- (38) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (Other name: 2C-E);
- (39) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (Other name: 2C-D);
- (40) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (Other name: 2C-C);
- (41) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (Other name: 2C-I);
- (42) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (Other name: 2C-T-2);
- (43) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (Other name: 2C-T-4);
- (44) 2-(2,5-Dimethoxyphenyl)ethanamine (Other name: 2C-H);
- (45) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (Other name: 2C-N);
- (46) 2-(2,5-Dimethoxy-4-(*n*)-propylphenyl)ethanamine (Other name: 2C-P);
- (47) 3,4-Methylenedioxy-*N*-methylcathinone (Other name: Methylone);
- (48) (1-Pentyl-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other names: UR-144, 1-pentyl-3-(2,2,3,3-tetramethylcyclopropyl)indole);
- (49) [1-(5-Fluoro-pentyl)-1*H*-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone (Other names: 5-fluoro-UR-144, 5-F-UR-144, XLR11, (5-fluoro-pentyl)-3-(2,2,3,3-tetramethylcyclopropyl)indole);
- (50) *N*-(1-Adamantyl)-1-pentyl-1*H*-indazole-3-carboxamide (Other names: APINACA, AKB48);
- (51) Quinolin-8-yl 1-pentyl-1*H*-indole-3-carboxylate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: PB-22; QUPIC);
- (52) Quinolin-8-yl 1-(5-fluoropentyl)-1*H*-indole-3-carboxylate, its optical, positional, and geometric isomers, salts and salts of isomers (Other names: 5-fluoro-PB-22; 5F-PB-22);
- (53) *N*-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers (Other name: AB-FUBINACA);
- (54) *N*-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1*H*-indazole-3-carboxamide (Other name: ADB-PINACA);
- (55) 2-(4-Iodo-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (Other names: 25I-NBOMe; 2CI-NBOMe; 25I; Cimbi-5);

- (56) 2-(4-Chloro-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82);
- (57) 2-(4-Bromo-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);
- *(58) Marijuana extract, meaning an extract containing one or more cannabinoids that has been derived from any plant of the genus *Cannabis*, except for extracts derived from hemp (as defined under Texas Agriculture Code 121) containing up to 0.3% delta-9-tetrahydrocannabinol on a dry weight basis, other than separated resin (whether crude or purified) obtained from the plant;
- (59) 4-Methyl-*N*-ethylcathinone (4-MEC);
- (60) 4-Methyl- α -pyrrolidinopropiophenone (4-MePPP);
- (61) α -Pyrrolidinopentiophenone ($[\alpha]$ -PVP);
- (62) 1-(1,3-Benzodioxol-5-yl)-2-(methylamino)butan-1-one (Other names: butylone; bk-MBDB);
- (63) 2-(Methylamino)-1-phenylpentan-1-one (Other name: pentedrone);
- (64) 1-(1,3-Benzodioxol-5-yl)-2-(methylamino)pentan-1-one (Other names: pentylone; bk-MBDP);
- (65) 4-Fluoro-*N*-methylcathinone (Other names: 4-FMC; flephedrone);
- (66) 3-Fluoro-*N*-methylcathinone (Other name: 3-FMC);
- (67) 1-(Naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one (Other name: naphyrone);
- (68) α -Pyrrolidinobutiophenone (Other name: $[\alpha]$ -PBP);
- (69) *N*-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1*H*-indazole-3-carboxamide (Other name: AB-CHMINACA);
- (70) *N*-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1*H*-indazole-3-carboxamide (Other name: AB-PINACA);
- (71) [1-(5-Fluoropentyl)-1*H*-indazol-3-yl](naphthalen-1-yl)methanone (Other name: THJ-2201);
- (72) 1-Methyl-4-phenyl-1,2,5,6-tetrahydro-pyridine (Other name: MPTP);
- (73) *N*-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1*H*-indazole-3-carboxamide (Other names: MAB-CHMINACA, ABD-CHMINACA);
- *(74) Methyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5F-ADB, 5F-MDMB-PINACA);
- *(75) Methyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3-methylbutanoate (Other name: 5F-AMB);
- *(76) *N*-(Adamantan-1-yl)-1-(5-fluoropentyl)-1*H*-indazole-3-carboxamide (Other names: 5F-APINACA; 5F-AKB48);
- *(77) *N*-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (Other name: ADB-FUBINACA);
- *(78) Methyl 2-(1-(cyclohexylmethyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (Other names: MDMB-CHMICA; MMB-CHMINACA);
- *(79) Methyl 2-(1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other name: MDMB-FUBINACA); and

*(80) Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (Other names: FUB-AMB; MMB-FUBINACA; AMB-FUBINACA).

-Schedule I depressants

Unless specifically excepted or unless listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following substances having a depressant effect on the central nervous system, including the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Gamma-hydroxybutyric acid (Other names: GHB; gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
- (2) Mecloqualone; and
- (3) Methaqualone.

-Schedule I stimulants

Unless specifically excepted or unless listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following substances having a stimulant effect on the central nervous system, including the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Aminorex (Other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine);
- (2) *N*-Benzylpiperazine (Other names: BZP; 1-benzylpiperazine), its optical isomers, salts and salts of isomers;
- (3) Cathinone (Other names: 2-amino-1-phenyl-1-propanone; α -aminopropiophenone; 2-aminopropiophenone; norephedrone);
- (4) Fenethylamine;
- (5) Methcathinone (Other names: 2-(methylamino)-propionophenone; α -(methylamino)propionophenone; 2-(methylamino)-1-phenylpropan-1-one; α -*N*-methylaminopropiophenone; monomethylpropion; ephedrone; *N*-methylcathinone; methylcathinone; AL-464; AL-422; AL-463; UR1432);
- (6) 4-Methylaminorex (Other names: U4Euh; McN-422);
- (7) *N*-Ethylamphetamine; and
- (8) *N,N*-Dimethylamphetamine (Other names: *N,N*- α -trimethylbenzene-ethaneamine; *N,N*- α -trimethylphenethylamine).

-Schedule I 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) The term 'cannabimimetic agents' 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:

(1-1) 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;

(1-2) 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;

(1-3) 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;

(1-4) 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; and,

(1-5) 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.

(2) 5-(1,1-Dimethylheptyl)-2-[(1*R*,3*S*)-3-hydroxycyclohexyl]-phenol (Other name: CP-47,497);

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

(4) 1-Pentyl-3-(1-naphthoyl)indole (Other names: JWH-018; AM678);

(5) 1-Butyl-3-(1-naphthoyl)indole (Other name: JWH-073);

(6) 1-Hexyl-3-(1-naphthoyl)indole (Other name: JWH-019);

(7) 1-[2-(4-Morpholinyl)ethyl]-3-(1-naphthoyl)indole (Other name: JWH-200);

(8) 1-Pentyl-3-(2-methoxyphenylacetyl)indole (Other name: JWH-250);

(9) 1-Pentyl-3-[1-(4-methoxynaphthoyl)]indole (Other name: JWH-081);

(10) 1-Pentyl-3-(4-methyl-1-naphthoyl)indole (Other name: JWH-122);

(11) 1-pentyl-3-(4-chloro-1-naphthoyl)indole (Other name: JWH-398);

(12) 1-(5-Fluoropentyl)-3-(1-naphthoyl)indole (Other name: AM2201);

(13) 1-(5-Fluoropentyl)-3-(2-iodobenzoyl)indole (Other name: AM694);

(14) 1-Pentyl-3-[(4-methoxy)-benzoyl]indole (Other names: SR-19, RCS-4);

(15) 1-Cyclohexylethyl-3-(2-methoxyphenylacetyl)indole (Other names: SR-18;RCS-8); and

(16) 1-Pentyl-3-(2-chlorophenylacetyl)indole (Other name: JWH-203).

-Schedule I temporarily listed substances subject to emergency scheduling by the U.S. Drug Enforcement Administration.

Unless specifically excepted or unless listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following substances or that contains any of the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

*(1) *N*-(1-Phenethylpiperidin-4-yl)-*N*-phenylpentanamide (Other name: valeryl fentanyl);

*(2) *N*-(4-Methoxyphenyl)-*N*-(1-phenethylpiperidin-4-yl)butyramide (Other name: *p*-methoxybutyryl fentanyl);

*(3) *N*-(4-Chlorophenyl)-*N*-(1-phenethylpiperidin-4-yl)isobutyramide (Other name: *p*-chloroisobutyryl fentanyl);

*(4) *N*-(1-Phenethylpiperidin-4-yl)-*N*-phenylisobutyramide (Other name: isobutyryl fentanyl);

*(5) *N*-(1-phenethylpiperidin-4-yl)-*N*-phenylcyclopentanecarboxamide (Other name: cyclopentyl fentanyl);

(6) Fentanyl-related substances.

(6-1) 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:

(6-1-1) Replacement of the phenyl portion of the phenethyl group by any monocycle, whether or not further substituted in or on the monocycle;

(6-1-2) Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxy, hydroxyl, halo, haloalkyl, amino or nitro groups;

(6-1-3) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxy, ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;

(6-1-4) Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the aromatic monocycle; and/or

(6-1-5) Replacement of the *N*-propionyl group by another acyl group.

(6-2) This definition includes, but is not limited to, the following substances:

(6-2-1) *N*-(1-(2-Fluorophenethyl)piperidin-4-yl)-*N*-(2-fluorophenyl)propionamide (Other name: 2'-fluoro-*o*-fluorofentanyl);

(6-2-2) *N*-(2-Methylphenyl)-*N*-(1-phenethylpiperidin-4-yl)acetamide (Other name: *o*-methyl acetylfentanyl);

(6-2-3) *N*-(1-Phenethylpiperidin-4-yl)-*N*,3-diphenylpropanamide (Other names: β' -phenyl fentanyl; hydrocinnamoyl fentanyl); and,

(6-2-4) *N*-(1-Phenethylpiperidin-4-yl)-*N*-phenylthiophene-2-carboxamide (Other name: thiofuranyl fentanyl).

*(7) Naphthalen-1-yl-1-(5-fluoropentyl)-1*H*-indole-3-carboxylate (Other names: NM2201; CBL2201);

*(8) *N*-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1*H*-indazole-3-carboxamide (Other name: 5F-AB-PINACA);

*(9) 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);

*(10) Methyl 2-(1-(cyclohexylmethyl)-1*H*-indole-3-carboxamido)-3-methylbutanoate (Other names: MMB-CHMICA; AMB-CHMICA);

*(11) 1-(5-Fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-pyrrolo[2,3-*b*]pyridine-3-carboxamide (Other name: 5F-CUMYL-P7AICA);

*(12) *N*-ethylpentylone (Other names: ephylone, 1-(1,3-benzodioxil-5-yl)-2-(ethylamino)-pentan-1-one);

(13) Ethyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other name: 5F-EDMB-PINACA);

(14) Methyl 2-(1-(5-fluoropentyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (Other name: 5F-MDMB-PICA);

(15) *N*-(Adamantan-1-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA; AKB48 *N*-(4-FLUOROBENZYL));

(16) 1-(5-Fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (Other names: 5F-CUMYL-PINACA; SGT-25);

(17) (1-(4-Fluorobenzyl)-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144);

(18) *N*-Ethylhexedrone (Other name: 2-(ethylamino)-1-phenylhexan-1-one);

(19) α -pyrrolidinohexanophenone (Other names: α -PHP; α -pyrrolidinohexiophenone; 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one);

(20) 4-Methyl- α -ethylaminopentiophenone (Other names: 4-MEAP; 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one);

(21) 4-Methyl- α -pyrrolidinohexiophenone (Other names: MPHP, 4'-methyl- α -pyrrolidinohexanophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);

(22) α -pyrrolidinoheptaphenone (Other names: PV8; 1-phenyl-2-(pyrrolidin-1-yl)heptan-1-one); and

(23) 4-Chloro- α -pyrrolidinovalerophenone (Other names: 4-chloro- α -PVP; 4-chloro- α -pyrrolidinopentiophenone; 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one); and,

*(24) *N,N*-diethyl-2-(2-(4 isopropoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts and salts of isomers, esters and ethers (Other names: isotonitazene; *N,N*-diethyl-2-[[4-(1-methylethoxy)phenyl]methyl]-5-nitro-1*H*-benzimidazole-1-ethanamine)

SCHEDULE II

Schedule II consists of:

Schedule II substances, vegetable origin or chemical synthesis

The following substances, however produced, except those narcotic drugs listed in other schedules:

(1) Opium and opiate, and a salt, compound, derivative, or preparation of opium or opiate, other than thebaine-derived butorphanol, naldemedine, naloxegol, naloxone and its salts, *6 β -naltrexol, naltrexone and its salts, and nalmeffene and its salts, but including:

- (1-1) Codeine (Other names: Morphine methyl ester; methyl morphine);
- (1-2) Dihydroetorphine (Other name: DHE);
- (1-3) Ethylmorphine (Other name: Dionin);
- (1-4) Etorphine hydrochloride (Other names: Etorphine HCl; M99);
- (1-5) Granulated opium;
- (1-6) Hydrocodone (Other name: dihydrocodeinone);
- (1-7) Hydromorphone (Other names: Dilaudid; dihydromorphinone);
- (1-8) Metopon;
- (1-9) Morphine (Other names: MS Contin; Roxanol; Oramorph; RMS; MSIR);
- (1-10) Noroxymorphone;
- (1-11) Opium extracts;
- (1-12) Opium fluid extracts;
- (1-13) Oripavine;
- (1-14) Oxycodone (Other names: OxyContin; Percocet; Endocet; Roxicodone; Roxicet);
- (1-15) Oxymorphone (Other name: Numorphan);
- (1-16) Powdered opium;
- (1-17) Raw opium (Other name: gum opium);
- (1-18) Thebaine; and
- (1-19) Tincture of opium (Other name: Laudanum).

(2) A salt, compound, isomer, derivative, or preparation of a substance that is chemically equivalent or identical to a substance described by paragraph (1) of Schedule II substances, vegetable origin or chemical synthesis, other than the isoquinoline alkaloids of opium;

(3) Opium poppy and poppy straw;

(4) Cocaine, including:

(4-1) its salts, its optical, position, and geometric isomers, and the salts of those isomers;

(4-2) Coca leaves and any salt, compound, derivative, or preparation of coca leaves and ecgonine and their salts, isomers, derivatives and salts of isomers and derivatives and any salt, compound derivative or preparation thereof which is chemically equivalent or identical to a substance described by this paragraph, except that the substances shall not include:

(4-2-1) Decocainized coca leaves or extractions of coca leaves which extractions do not that do not contain cocaine or ecgonine; or

(4-2-2) Ioflupane.

(5) Concentrate of poppy straw, meaning the crude extract of poppy straw in liquid, solid, or powder form that contains the phenanthrene alkaloids of the opium poppy.

Schedule II opiates

The following opiates, including their isomers, esters, ethers, salts, and salts of isomers, if the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

- (1) Alfentanil (Other name: Alfenta);
- (2) Alphaprodine (Other name: Nisentil);
- (3) Anileridine (Other name: Leritine);
- (4) Bezitramide (Other name: Burgodin);
- (5) Carfentanil (Other name: Wildnil);
- (6) Dextropropoxyphene, bulk (nondosage form) (Other name: Propoxyphene);
- (7) Dihydrocodeine (Other names: Didrate; Parzone);
- (8) Diphenoxylate;
- (9) Fentanyl (Other names: Duragesic; Oralet; Actiq; Sublimaze; Innovar);
- (10) Isomethadone (Other name: Isoamidone);
- (11) Levo-alpha-acetylmethadol (Other names: levo- α -acetylmethadol; levomethadyl acetate, LAAM);
- (12) Levomethorphan;
- (13) Levorphanol;
- (14) Metazocine;
- (15) Methadone (Other names: Dolophine; Methadose; Amidone);
- (16) Methadone Intermediate (Other name: 4-cyano-2-dimethylamino-4,4-diphenyl butane);
- (17) Moramide Intermediate (Other name: 2-methyl-3-morpholino-1,1-diphenylpropane-carboxylic acid);
- *(18) Oliceridine (Other name: N-[(3-methoxythiophen-2-yl)methyl] ({2-[(9R)-9-(pyridin-2-yl)-6-oxaspiro [4.5]decan-9-yl]ethyl})amine fumarate)
- (19) Pethidine (meperidine);
- (20) Pethidine Intermediate-A (Other name: 4-cyano-1-methyl-4-phenylpiperidine);
- (21) Pethidine Intermediate-B (Other name: ethyl-4-phenylpiperidine-4-carboxylate);
- (22) Pethidine Intermediate-C (Other name: 1-methyl-4-phenylpiperidine-4-carboxylic acid);
- (23) Phenazocine (Other names: Narphen; Prinadol);
- (24) Piminodine;
- (25) Racemethorphan;
- (26) Racemorphan (Other name: Dromoran);
- (27) Remifentanil (Other name: Ultiva);
- (28) Sufentanil (Other name: Sufenta);
- (29) Tapentadol; and
- (30) Thiafentanil (Other name: methyl 4-(2-methoxy-N-phenylacetamido)-1-(2-(thiophen-2-yl)ethyl)piperidine-4-carboxylate).

• Schedule II stimulants

Unless listed in another schedule and except as provided by the Texas Controlled Substances Act, Health and Safety Code, Section 481.033, a material, compound, mixture, or preparation that contains any quantity of the following substances having a potential for abuse associated with a stimulant effect on the central nervous system:

- (1) Amphetamine, its salts, optical isomers, and salts of its optical isomers;
- (2) Methamphetamine, including its salts, optical isomers, and salts of optical isomers;
- (3) Methylphenidate and its salts;
- (4) Phenmetrazine and its salts; and
- (5) Lisdexamfetamine, including its salts, isomers, and salts of its isomers.

-Schedule II depressants

Unless listed in another schedule, a material, compound, mixture or preparation that contains any quantity of the following substances having a depressant effect on the central nervous system, including the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Amobarbital;
- (2) Glutethimide;
- (3) Pentobarbital; and
- (4) Secobarbital.

-Schedule II hallucinogenic substances

(1) Nabilone (Another name for nabilone: (\pm)-*trans*-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1-hydroxy-6,6-dimethyl-9*H*-dibenzo[b,d]pyran-9-one); and

(2) Dronabinol in oral solution in drug products approved for marketing by the U.S. Food and Drug Administration.

-Schedule II precursors

Unless specifically excepted or listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following substances:

- (1) Immediate precursor to methamphetamine:
 - (1-1) Phenylacetone and methylamine if possessed together with intent to manufacture methamphetamine;
- (2) Immediate precursor to amphetamine and methamphetamine:
 - (2-1) Phenylacetone (Other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone);
- (3) Immediate precursors to phencyclidine (PCP):
 - (3-1) 1-phenylcyclohexylamine;
 - (3-2) 1-piperidinocyclohexanecarbonitrile (PCC); and

- (4) Immediate precursor to fentanyl:
(4-1) 4-anilino-N-phenethylpiperidine (ANPP).

SCHEDULE III

Schedule III consists of:

-Schedule III depressants

Unless listed in another schedule and except as provided by the Texas Controlled Substances Act, Health and Safety Code, Section 481.033, a material, compound, mixture, or preparation that contains any quantity of the following substances having a potential for abuse associated with a depressant effect on the central nervous system:

- (1) A compound, mixture, or preparation containing amobarbital, secobarbital, pentobarbital, or any of their salts and one or more active medicinal ingredients that are not listed in a schedule;
- (2) A suppository dosage form containing amobarbital, secobarbital, pentobarbital, or any of their salts and approved by the U.S. Food and Drug Administration for marketing only as a suppository;
- (3) A substance that contains any quantity of a derivative of barbituric acid, or any salt of a derivative of barbituric acid, except those substances that are specifically listed in other schedules;
- (4) Chlorhexadol;
- (5) Any drug product containing gamma hydroxybutyric acid, including its salts, isomers, and salts of isomers, for which an application is approved under Section 505 of the Federal Food Drug and Cosmetic Act.
- (6) Ketamine, its salts, isomers, and salts of isomers. Some other names for ketamine: (\pm)-2-(2-chlorophenyl)-2-(methylamino)-cyclohexanone;
- (7) Lysergic acid;
- (8) Lysergic acid amide;
- (9) Methyprylon;
- (10) Perampanel, and its salts, isomers, and salts of isomers;
- (11) Sulfondiethylmethane;
- (12) Sulfonethylmethane;
- (13) Sulfonmethane; and,
- (14) Tiletamine and zolazepam or any salt thereof. (Some trade or other names for a tiletamine zolazepam combination product: Telazol. Some trade or other names for tiletamine: 2-(ethylamino)-2-(2-thienyl) cyclohexanone. Some trade or other names for zolazepam: 4-(2-fluorophenyl)-6,8-dihydro-1,3,8 trimethyl-pyrazolo[3,4-e][1,4] diazepin-7(1H)-one, flupyrzapon.)

-Schedule III

- (1) Nalorphine.

-Schedule III narcotics

Unless specifically excepted or unless listed in another schedule:

(1) A material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs, or any of their salts:

(1-1) not more than 1.8 grams of codeine, or any of its salts, per 100 milliliters or not more than 90 milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium;

(1-2) not more than 1.8 grams of codeine, or any of its salts, per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(1-3) not more than 1.8 grams of dihydrocodeine, or any of its salts, per 100 milliliters or not more than 90 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts;

(1-4) not more than 300 milligrams of ethylmorphine, or any of its salts, per 100 milliliters or not more than 15 milligrams per dosage unit, with one or more active, non-narcotic ingredients in recognized therapeutic amounts;

(1-5) not more than 500 milligrams of opium per 100 milliliters or per 100 grams, or not more than 25 milligrams per dosage unit, with one or more active, nonnarcotic ingredients in recognized therapeutic amounts; and

(1-6) not more than 50 milligrams of morphine, or any of its salts, per 100 milliliters or per 100 grams with one or more active, nonnarcotic ingredients in recognized therapeutic amounts; and

(2) any material, compound, mixture, or preparation containing any of the following narcotic drugs or their salts:

(2-1) Buprenorphine.

-Schedule III stimulants

Unless listed in another schedule, a material, compound, mixture or preparation that contains any quantity of the following substances having a stimulant effect on the central nervous system, including the substance's salts, optical, position, or geometric isomers, and salts of the substance's isomers, if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Benzphetamine;
- (2) Chlorphentermine;
- (3) Clortermine; and
- (4) Phendimetrazine.

-Schedule III anabolic steroids and hormones

Anabolic steroids, including any drug or hormonal substance, chemically and pharmacologically related to testosterone (other than estrogens, progestins, corticosteroids, and dehydroepiandrosterone), and include the following:

- (1) Androstenediol--
 - (1-1) 3 β ,17 β -dihydroxy-5 α -androstane;
 - (1-2) 3 α ,17 β -dihydroxy-5 α -androstane;
- (2) Androstenedione (5 α -androstane-3,17-dione);
- (3) Androstenediol--
 - (3-1) 1-androstenediol (3 β ,17 β -dihydroxy-5 α -androst-1-ene);
 - (3-2) 1-androstenediol (3 α ,17 β -dihydroxy-5 α -androst-1-ene);
 - (3-3) 4-androstenediol (3 β ,17 β -dihydroxy-androst-4-ene);
 - (3-4) 5-androstenediol (3 β ,17 β -dihydroxy-androst-5-ene);
- (4) Androstenedione--
 - (4-1) 1-androstenedione (5 α -androst-1-en-3,17-dione);
 - (4-2) 4-androstenedione (androst-4-en-3,17-dione);
 - (4-3) 5-androstenedione (androst-5-en-3,17-dione);
- (5) Bolasterone (7 α ,17 α -dimethyl-17 β -hydroxyandrost-4-en-3-one);
- (6) Boldenone (17 β -hydroxyandrost-1,4,-diene-3-one);
- (7) Boldione (androst-1,4-diene-3,17-dione);
- (8) Calusterone (7 β ,17 α -dimethyl-17 β -hydroxyandrost-4-en-3-one);
- (9) Clostebol (4-chloro-17 β -hydroxyandrost-4-en-3-one);
- (10) Dehydrochloromethyltestosterone (4-chloro-17 β -hydroxy-17 α -methyl-androst-1,4-dien-3-one);
- (11) Δ -1-Dihydrotestosterone (17 β -hydroxy-5 α -androst-1-en-3-one)
(Other Name: 1-testosterone);
- (12) Desoxymethyltestosterone (17 α -methyl-5 α -androst-2-en-17 β -ol; madol);
- (13) 4-Dihydrotestosterone (17 β -hydroxy-androstan-3-one);
- (14) Drostanolone (17 β -hydroxy-2 α -methyl-5 α -androstane-3-one);
- (15) Ethylestrenol (17 α -ethyl-17 β -hydroxyestr-4-ene);
- (16) Fluoxymesterone (9-fluoro-17 α -methyl-11 β ,17 β -dihydroxyandrost-4-en-3-one);
- (17) Formebolone (2-formyl-17 α -methyl-11 α ,17 β -dihydroxyandrost-1,4-dien-3-one);
- (18) Furazabol (17 α -methyl-17 β -hydroxyandrostano[2,3-c]-furazan);
- (19) 13 β -Ethyl-17 β -hydroxygon-4-en-3-one;
- (20) 4-Hydroxytestosterone (4,17 β -dihydroxy-androst-4-en-3-one);
- (21) 4-Hydroxy-19-nortestosterone (4,17 β -dihydroxy-estr-4-en-3-one);
- (22) Mestanolone (17 α -methyl-17 β -hydroxy-5 α -androstane-3-one);
- (23) Mesterolone (1 α -methyl-17 β -hydroxy-5 α -androstane-3-one);
- (24) Methandienone (17 α -methyl-17 β -hydroxyandrost-1,4-dien-3-one);
- (25) Methandriol (17 α -methyl-3 β ,17 β -dihydroxyandrost-5-ene);
- (26) Methenolone (1-methyl-17 β -hydroxy-5 α -androst-1-en-3-one);

- (27) 17 α -Methyl-3 β , 17 β -dihydroxy-5 α -androstane;
- (28) Methasterone (2 α ,17 α -dimethyl-5 α -androst-17 β -ol-3-one);
- (29) 17 α -Methyl-3 α ,17 β -dihydroxy-5 α -androstane;
- (30) 17 α -Methyl-3 β ,17 β -dihydroxyandrost-4-ene;
- (31) 17 α -Methyl-4-hydroxynandrolone (17 α -methyl-4-hydroxy-17 β -hydroxyestr-4-en-3-one);
- (32) Methyldienolone (17 α -methyl-17 β -hydroxyestra-4,9(10)-dien-3-one);
- (33) 17 α -Methyl-17 β -hydroxyestra-4,9-11-trien-3-one;
- (34) Methyltestosterone (17 α -methyl-17 β -hydroxyandrost-4-en-3-one);
- (35) Mibolerone (7 α ,17 α -dimethyl-17 β -hydroxyestr-4-en-3-one);
- (36) 17 α -Methyl-delta-1-dihydrotestosterone (17 β -hydroxy-17 α -methyl-5 α -androst-1-en-3-one) (Other name: 17 α -methyl-1-testosterone);
- (37) Nandrolone (17 β -hydroxyestr-4-en-3-one);
- (38) Norandrostenediol--
 - (38-1) 19-Nor-4-androstenediol (3 β ,17 β -dihydroxyestr-4-ene);
 - (38-2) 19-Nor-4-androstenediol (3 α ,17 β -dihydroxyestr-4-ene);
 - (38-3) 19-Nor-5-androstenediol (3 β ,17 β -dihydroxyestr-5-ene);
 - (38-4) 19-Nor-5-androstenediol (3 α ,17 β -dihydroxyestr-5-ene);
- (39) Norandrostenedione--
 - (39-1) 19-Nor-4-androstenedione (estr-4-en-3,17-dione);
 - (39-2) 19-Nor-5-androstenedione (estr-5-en-3,17-dione);
- (40) 19-Nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione);
- (41) Norbolethone (13 β ,17 α -diethyl-17 β -hydroxygon-4-en-3-one);
- (42) Norclostebol (4-chloro-17 β -hydroxyestr-4-en-3-one);
- (43) Norethandrolone (17 α -ethyl-17 β -hydroxyestr-4-en-3-one);
- (44) Normethandrolone (17 α -methyl-17 β -hydroxyestr-4-en-3-one);
- (45) Oxandrolone (17 α -methyl-17 β -hydroxy-2-oxa-5 α -androst-3-one);
- (46) Oxymesterone (17 α -methyl-4,17 β -dihydroxyandrost-4-en-3-one);
- (47) Oxymetholone (17 α -methyl-2-hydroxymethylene-17 β -hydroxy-5 α -androst-3-one);
- (48) Prostanazol (17 β -hydroxy-5 α -androstano[3,2-c]pyrazole);
- (49) Stanozolol (17 α -methyl-17 β -hydroxy-5 α -androst-2-eno[3,2-c]-pyrazole);
- (50) Stenbolone (17 β -hydroxy-2-methyl-5 α -androst-1-en-3-one);
- (51) Testolactone (13-hydroxy-3-oxo-13,17-secoandrost-1,4-dien-17-oic acid lactone);
- (52) Testosterone (17 β -hydroxyandrost-4-en-3-one);
- (53) Tetrahydrogestrinone (13 β ,17 α -diethyl-17 β -hydroxygon-4,9,11-trien-3-one);
- (54) Trenbolone (17 β -hydroxyestr-4,9,11-trien-3-one); and
- (55) any salt, ester, or ether of a drug or substance.

-Schedule III hallucinogenic substances

(1) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in U.S. Food and Drug Administration approved drug product.

(Some other names for dronabinol:

(6a*R*-*trans*)-6a,7,8,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6*H*-dibenzo[b,d]pyran-1-ol; (-)- Δ -9-(*trans*)-tetrahydrocannabinol).

SCHEDULE IV

Schedule IV consists of:

-Schedule IV depressants

Except as provided by the Texas Controlled Substances Act, Health and Safety Code, Section 481.033, a material, compound, mixture, or preparation that contains any quantity of the following substances or any of the substance's salts, isomers, and salts of isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation having a potential for abuse associated with a depressant effect on the central nervous system:

- (1) Alfaxalone (5 α -pregnan-3 α -ol-11,20-dione);
- (2) Alprazolam;
- (3) Barbital;
- (4) Brexanolone (3 α -hydroxy-5 α -pregnan-20-one) (Other name: allopregnanolone);
- (5) Bromazepam;
- (6) Camazepam;
- (7) Chloral betaine;
- (8) Chloral hydrate;
- (9) Chlordiazepoxide;
- (10) Clobazam;
- (11) Clonazepam;
- (12) Clorazepate;
- (13) Clotiazepam;
- (14) Cloxazolam;
- (15) Delorazepam;
- (16) Diazepam;
- (17) Dichloralphenazone;
- (18) Estazolam;
- (19) Ethchlorvynol;
- (20) Ethinamate;
- (21) Ethyl loflazepate;
- (22) Fludiazepam;
- (23) Flunitrazepam;
- (24) Flurazepam;
- (25) Fospropofol;

- (26) Halazepam;
- (27) Haloxazolam;
- (28) Ketazolam;
- *(29) Lemborexant;
- (30) Loprazolam;
- (31) Lorazepam;
- (32) Lormetazepam;
- (33) Mebutamate;
- (34) Medazepam;
- (35) Meprobamate;
- (36) Methohexital;
- (37) Methylphenobarbital (mephobarbital);
- (38) Midazolam;
- (39) Nimetazepam;
- (40) Nitrazepam;
- (41) Nordiazepam;
- (42) Oxazepam;
- (43) Oxazolam;
- (44) Paraldehyde;
- (45) Petrichloral;
- (46) Phenobarbital;
- (47) Pinazepam;
- (48) Prazepam;
- (49) Quazepam;
- *(50) Remimazolam;
- (51) Suvorexant;
- (52) Temazepam;
- (53) Tetrazepam;
- (54) Triazolam;
- (55) Zaleplon;
- (56) Zolpidem; and
- (57) Zopiclone.

-Schedule IV stimulants

Unless listed in another schedule, a material, compound, mixture, or preparation that contains any quantity of the following substances having a stimulant effect on the central nervous system, including the substance's salts, optical, position, or geometric isomers, and salts of those isomers if the existence of the salts, isomers, and salts of isomers is possible within the specific chemical designation:

- (1) Cathine [(+)-norpseudoephedrine];
- (2) Diethylpropion;
- (3) Fencamfamin;
- (4) Fenfluramine;
- (5) Fenproporex;
- (6) Mazindol;
- (7) Mefenorex;
- (8) Modafinil;

- (9) Pemoline (including organometallic complexes and their chelates);
- (10) Phentermine;
- (11) Pipradrol;
- (12) Sibutramine;
- (13) Solriamfetol ((*R*)-2-amino-3-phenylpropyl carbamate) (Other names: benzenepropanol; β -amino-carbamate (ester));
and
- (14) SPA [1-dimethylamino-1,2-diphenylethane].

-Schedule IV narcotics

Unless specifically excepted or unless listed in another schedule, a material, compound, mixture, or preparation containing limited quantities of the following narcotic drugs or their salts:

- (1) Not more than 1 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit;
- (2) Dextropropoxyphene (α -(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane); and
- (3) 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol (other name: tramadol).

-Schedule IV other substances

Unless specifically excepted or unless listed in another schedule, a material, compound, substance's salts:

- (1) Butorphanol, including its optical isomers;
- (2) Carisoprodol;
- (3) Eluxadolone (other name: 5-[[[(2*S*-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-oxopropyl][(1*S*)-1-(4-phenyl-1*H*-imidazol-2-yl)ethyl]amino]methyl]-2-methoxybenzoic acid) including its salts, isomers, and salts of isomers;
- (4) Lorcarserin including its salts, isomers and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers is possible; and
- (5) Pentazocine, its salts, derivatives, compounds, or mixtures.

SCHEDULE V

Schedule V consists of:

-Schedule V narcotics containing non-narcotic active medicinal ingredients

A compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs that also contain one or more nonnarcotic active medicinal ingredients in sufficient proportion to confer on the compound, mixture or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:

- (1) Not more than 200 milligrams of codeine, or any of its salts, per 100 milliliters or per 100grams;
- (2) Not more than 100 milligrams of dihydrocodeine, or any of its salts, per 100 milliliters or per 100 grams;
- (3) Not more than 100 milligrams of ethylmorphine, or any of its salts, per 100 milliliters or per 100 grams;
- (4) Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of atropine sulfate per dosage unit;
- (5) Not more than 15 milligrams of opium per 29.5729 milliliters or per 28.35 grams; and
- (6) Not more than 0.5 milligram of difenoxin and not less than 25 micrograms of atropine sulfate per dosage unit.

-Schedule V stimulants

Unless specifically exempted or excluded or unless listed in another schedule, a 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) Pyrovalerone.

-Schedule V depressants

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

- (1) Brivaracetam ((2*S*)-2-[(4*R*)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (Other names; BRV; UCB-34714; Briviact);
- * (2) Cenobamate [(1*R*-1-(2-chlorophenyl)-2-(tetrazol-2-yl)ethyl] carbamate;
- (3) Ezogabine including its salts, isomers and salts of isomers, whenever the existence of such salts, isomers and salts of isomers is possible;
- (4) Lacosamide [(*R*)-2-acetoamido-*N*-benzyl-3-methoxy-propionamide];
- * (5) Lasmiditan [2,4,6-trifluoro-*N*-(6-(1-methylpiperidine-4-carbonyl)pyridine-2-yl)-benzamide]; and
- (6) Pregabalin [(*S*)-3-(aminomethyl)-5-methylhexanoic acid].