

IMPOUND ORDERS ISSUED:

Name	Type of Order	License Number	Address	Action	Date of Issuance
Texas Gamma Ray, LLC DBA TGR Industrial Services	Impound Order	L05561 (Site 005)	8777 Tallyho Road, Building 1 Houston, Texas	Impound industrial radiography devices	06/13/22
Texas Gamma Ray, LLC DBA TGR Industrial Services	Impound Order	L05561 (Site 006)	7080 Mayard Road Houston, Texas	Impound industrial radiography device	06/13/22

TRD-202202818
 Cynthia Hernandez
 General Counsel
 Department of State Health Services
 Filed: July 27, 2022

Order Adding Daridorexant to Schedule IV, Permanently Placing Five Synthetic Cannabinoids in Schedule I, and Temporarily Placing Seven Synthetic Benzimidazole-opioid Substances in Schedule I



The Drug Enforcement Administration (DEA) issued an interim final rule placing daridorexant in schedule IV of the Controlled Substances Act, including its salts, isomers of salts, and salts of isomers whenever the existence of such salts, isomers, and salts of such isomers is possible within the specific chemical designation. The interim final rule was published in the April 7, 2022 edition of the *Federal Register*, Volume 87, Number 67, pages 20313-20318 and was effective April 7, 2022. This action was taken based on the following:

- (1) Daridorexant has a low potential for abuse relative to the drugs or other substances in schedule III;
- (2) Daridorexant has a currently accepted medical use in treatment in the United States; and
- (3) Abuse of daridorexant may lead to limited physical dependence or psychological dependence relative to the drugs or other substances in schedule III.

The DEA issued a final rule permanently placing the following substances in schedule I: ethyl 2-(1-(5-fluoropentyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate (Other name: 5F-EDMB-PINACA); methyl 2-(1-(5-fluoropentyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5F-MDMB-PICA; 5F-MDMB-2201); *N*-(adamantan-1-yl)-1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamide (Other names: FUB-AKB48; FUB-APINACA, AKB48 *N*-(4-FLUOROBENZYL)); 1-(5-fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (Other names: 5F-CUMYL-PINACA; SGT-25); and (1-(4-fluorobenzyl)-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144) including 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. The final rule was published in the April 7, 2022 edition of the *Federal Register*, Volume 87, Number 67, pages 20318-20321 and was effective April 7, 2022. This action was taken based on the following:

- (1) 5F-EDMB-PINACA, 5F-MDMB-PICA, FUB-AKB48, 5F-CUMYL-PINACA, and

FUB-144 have a high potential for abuse that is comparable to other schedule I substances such as JWH-018;

(2) 5F-EDMB-PINACA, 5F-MDMB-PICA, FUB-AKB48, 5F-CUMYL-PINACA, and FUB-144 currently have no accepted medical use in treatment in the United States; and,

(3) There is a lack of accepted safety for use of 5F-EDMB-PINACA, 5F-MDMB-PICA, FUB-AKB48, 5F-CUMYL-PINACA, and FUB-144 under medical supervision.

The DEA issued a temporary amendment temporarily placing the following seven synthetic benzimidazole-opioid substances into schedule I: 2-(2-(4-butoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine (Other name: butonitazene); 2-(2-(4-ethoxybenzyl)-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine (Other names: etodesnitazene; etazene); *N,N*-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: flunitazene); *N,N*-diethyl-2-(2-(4-methoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: metodesnitazene); *N,N*-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: metonitazene); 2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (Other names: *N*-pyrrolidino etonitazene; etonitazepyne); and *N,N*-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: protonitazene) including 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. The temporary amendment was published in the April 12, 2022 edition of the *Federal Register*, Volume 87, Number 70, pages 21556-21561. This action was taken based on the following:

1. Temporary placement of butonitazene, etodesnitazene, flunitazene, metodesnitazene, metonitazene, *N*-pyrrolidino etonitazene and protonitazene into schedule I is necessary to avoid an imminent hazard to public safety;
2. Butonitazene, etodesnitazene, flunitazene, metodesnitazene, metonitazene, *N*-pyrrolidino etonitazene and protonitazene have a high potential for abuse;

3. Butonitazene, etodesnitazene, flunitazene, metodesnitazene, metonitazene, *N*-pyrrolidino etonitazene and protonitazene have no currently accepted medical use in the United States; and,

4. There is a lack of accepted safety for use of butonitazene, etodesnitazene, flunitazene, metodesnitazene, metonitazene, *N*-pyrrolidino etonitazene and protonitazene under medical supervision.

Pursuant to Section 481.034(g), as amended by the 75th legislature, of the Texas Controlled Substances Act, Health and Safety Code, Chapter 481, at least thirty-one days have expired since notice of the above referenced actions were published in the Federal Register. In the capacity as Commissioner of the Texas Department of State Health Services, John Hellerstedt, M.D., does hereby order that the substance daridorexant be added to schedule IV; the substances 5F-EDMB-PINACA, 5F-MDMB-PICA, FUB-AKB48, 5F-CUMYL-PINACA, and FUB-144 be permanently placed into schedule I; and the substances butonitazene, etodesnitazene, flunitazene, metodesnitazene, metonitazene, *N*-pyrrolidino etonitazene, and protonitazene be temporarily placed into schedule I.

-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 name: MDMA);
- (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 name: *N*-hydroxy MDA);
- (14) 3,4,5-Trimethoxyamphetamine (Other name: 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 tetrahydrocannabinols in hemp (as defined under Section 297A(1) of the Agricultural Marketing Act of 1946), 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 phenylcyclohexyl)-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*-methylethylcathinone; mephedrone);

(37) 3,4-Methylenedioxypropylvalerone (Other name: 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: Methydone);
- (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-Fluoropentyl)-1*H*-indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone (Other names: 5-fluoro-UR-144; 5-F-UR-144, XLR11; (5-fluoropentyl)-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*, other than separated resin (whether crude or purified) obtained from the plant;
- (59) 4-Methyl-*N*-ethylcathinone (Other name: 4-MEC);
- (60) 4-Methyl- α -pyrrolidinopropiophenone (Other name: 4-MePPP);
- (61) α -Pyrrolidinopentiophenone (Other name: [α]-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: α -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);
(80) Methyl 2-(1-(4-fluorobenzyl)-1*H*-indazole-3-carboxamido)-3-methylbutanoate (Other names: FUB-AMB; MMB-FUBINACA; AMB-FUBINACA);
(81) Naphthalen-1-yl-1-(5-fluoropentyl)-1*H*-indole-3-carboxylate (Other names: NM2201; CBL2201);
(82) *N*-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1*H*-indazole-3-carboxamide (Other name: 5F-AB-PINACA);
(83) 1-(4-Cyanobutyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (Other names: 4-CN-CUMYL-BUTINACA; 4-cyano-CUMYL-BUTINACA; 4-CN-CUMYL-BINACA; CUMYL-4CN-BINACA; SGT-78);
(84) Methyl 2-(1-(Cyclohexylmethyl)-1*H*-indole-3-carboxamido)-3-methylbutanoate (Other names: MMB-CHMICA; AMB-CHMICA);
(85) 1-(5-Fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-pyrrolo[2,3-*b*]pyridine-3-carboxamide (Other name: 5F-CUMYL-P7AICA);

(86) 1-(1,3-Benzodioxol-5-yl)-2-(ethylamino)pentan-1-one (Other names: *N*-ethylpentylone; ephylone);

(87) Methyl 2-(1-(4-fluorobutyl)-1*H*-indazole-3-carboxamido)-3,3-dimethylbutanoate) (Other names: 4F-MDMB-BINACA; 4F-MDMB-BUTINACA);

(88) 1-(4-Methoxyphenyl)-*N*-methylpropan-2-amine (Other names: *p*-methoxymethamphetamine; PMMA);

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

*(90) Methyl 2-(1-(5-fluoropentyl)-1*H*-indole-3-carboxamido)-3,3-dimethylbutanoate (Other names: 5F-MDMB-PICA; 5F-MDMB-2201);

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

*(92) 1-(5-Fluoropentyl)-*N*-(2-phenylpropan-2-yl)-1*H*-indazole-3-carboxamide (Other names: 5F-CUMYL-PINACA; SGT-25); and,

*(93) (1-(4-Fluorobenzyl)-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone (Other name: FUB-144).

-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 isomers, esters, ethers, salts and salts of isomers, esters, and ethers if the existence of the salts, esters, ethers isomers, and salts of isomers, esters, ethers is possible within the specific chemical designation:

(1) Fentanyl-related substances.

(1-1) 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:

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

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

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

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

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

(1-2) This definition includes, but is not limited to, the following substances:

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

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

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

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

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

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

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

(5) 4'-Methyl- α -pyrrolidinohexiophenone (Other names: MPHP; 4'-methyl- α -pyrrolidinohexanophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)hexan-1-one);

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

(7) 4'-Chloro- α -pyrrolidinovalerophenone (Other names: 4-chloro- α -PVP; 4'-chloro- α -pyrrolidinopentiophenone; 1-(4-chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one);

(8) 1-(1-(1-(4-Bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2*H*-benzo[d]imidazol-2-one (Other names: bromphine; 1-[1-[1-(4-bromophenyl)ethyl]-4-piperidinyl]-1,3-dihydro-2*H*-benzimidazol-2-one);

- * (9) 2-(2-(4-Butoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine (Other name: butonitazene);
- * (10) 2-(2-(4-Ethoxybenzyl)-1*H*-benzimidazol-1-yl)-*N,N*-diethylethan-1-amine (Other names: etodesnitazene; etazene);
- * (11) *N,N*-Diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: flunitazene);
- * (12) *N,N*-Diethyl-2-(2-(4-methoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: metodesnitazene);
- * (13) *N,N*-Diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: metonitazene);
- * (14) 2-(4-Ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (Other names: *N*-pyrrolidino etonitazene; etonitazepyne);
- and,
- * (15) *N,N*-Diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: protonitazene).

-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 (5α-pregnan-3α-ol-11,20-dione);
- (2) Alprazolam;
- (3) Barbitol;
- (4) Brexanolone (3α-hydroxy-5α-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) Daridorexant;
- (16) Delorazepam;
- (17) Diazepam;
- (18) Dichloralphenazone;
- (19) Estazolam;
- (20) Ethchlorvynol;
- (21) Ethinamate;
- (22) Ethyl loflazepate;
- (23) Fludiazepam;
- (24) Flunitrazepam;
- (25) Flurazepam;
- (26) Fospropofol;
- (27) Halazepam;
- (28) Haloxazolam;
- (29) Ketazolam;
- (30) Lemborexant;
- (31) Loprazolam;
- (32) Lorazepam;
- (33) Lormetazepam;
- (34) Mebutamate;
- (35) Medazepam;
- (36) Meprobamate;
- (37) Methohexital;
- (38) Methylphenobarbital (mephobarbital);
- (39) Midazolam;
- (40) Nimetazepam;
- (41) Nitrazepam;
- (42) Nordiazepam;
- (43) Oxazepam;
- (44) Oxazolam;
- (45) Paraldehyde;
- (46) Petrichloral;
- (47) Phenobarbital;
- (48) Pinazepam;
- (49) Prazepam;

- (50) Quazepam;
- (51) Remimazolam;
- (52) Suvorexant;
- (53) Temazepam;
- (54) Tetrazepam;
- (55) Triazolam;
- (56) Zaleplon;
- (57) Zolpidem; and
- (58) Zopiclone.

TRD-202202816
 Cynthia Hernandez
 General Counsel
 Department of State Health Services
 Filed: July 27, 2022



Texas Department of Insurance

Company Licensing

Application to do business in the state of Texas for AmFirst Specialty Insurance Company, a foreign life, accident and/or health company. The home office is in Ridgeland, Mississippi.

Application to do business in the state of Texas for CM Indemnity Insurance Company, a foreign fire and/or casualty company. The home office is in Merrill, Washington.

Application for United Benefit Life Insurance Company, a foreign life, accident and/or health company, to change its name to Cigna Insurance Company. The home office is in Cleveland, Ohio.

Any objections must be filed with the Texas Department of Insurance, within twenty (20) calendar days from the date of the *Texas Register* publication, addressed to the attention of John Carter, 333 Guadalupe Street, MC FRD-CL, Austin, Texas 78701.

TRD-202202827
 Justin Beam
 Chief Clerk
 Texas Department of Insurance
 Filed: July 27, 2022



Texas Lottery Commission

Scratch Ticket Game Number 2439 "\$5,000 Cash Blowout"

1.0 Name and Style of Scratch Ticket Game.

A. The name of Scratch Ticket Game No. 2439 is "\$5,000 CASH BLOWOUT". The play style is "key number match".

1.1 Price of Scratch Ticket Game.

A. The price for Scratch Ticket Game No. 2439 shall be \$1.00 per Scratch Ticket.

1.2 Definitions in Scratch Ticket Game No. 2439.

A. Display Printing - That area of the Scratch Ticket outside of the area where the overprint and Play Symbols appear.

B. Latex Overprint - The removable scratch-off covering over the Play Symbols on the front of the Scratch Ticket.

C. Play Symbol - The printed data under the latex on the front of the Scratch Ticket that is used to determine eligibility for a prize. Each Play Symbol is printed in Symbol font in black ink in positive except for dual-image games. The possible black Play Symbols are: 01, 03, 04, 05, 06, 07, 08, 09, 10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20, 2X SYMBOL, \$1.00, \$2.00, \$5.00, \$10.00, \$20.00, \$50.00, \$100 and \$5,000.

D. Play Symbol Caption - The printed material appearing below each Play Symbol which explains the Play Symbol. One caption appears under each Play Symbol and is printed in caption font in black ink in positive. The Play Symbol Caption which corresponds with and verifies each Play Symbol is as follows: