hearing, please submit an electronic copy in accessible PDF format that will be distributed to the appropriate DSHS staff. Handouts are limited to two pages (paper size: 8.5" by 11", one side only). Handouts must be emailed to DSHSPublicHearings@dshs.texas.gov immediately after pre-registering, but no later than 5:00 p.m., Sunday, April 28, 2024, and include the name of the person who will be commenting. Do not include health or other confidential information in your comments or handouts. Staff will not read handouts aloud during the public hearing, but handouts will be provided to the appropriate DSHS staff.

**Oral comments provided in-person at the public hearing location:** Members of the public may provide oral public comments during the public hearing in person at the public hearing location by pre-registering online using the Public Comment Registration form at https://www.surveymonkey.com/r/L8W7NF7. Do not include health information or other confidential information in your comments or handouts.

Written comments: A member of the public who wishes to provide written public comments only must email the comments to DSHSPublicHearings@dshs.texas.gov no later than 5:00 p.m., Sunday, April 28, 2024. Please include your name and the organization you are representing or that you are commenting as a private citizen. Written comments are limited to two pages (paper size: 8.5" by 11", one side only). Do not include health or other confidential information in your comments. Staff will not read written comments aloud during the public hearing, but comments will be provided to the appropriate DSHS staff.

Note: These procedures may be revised at the discretion of DSHS.

**Contact:** Questions regarding agenda items, content, or public hearing arrangements should be directed to:

Mail: Department of State Health Services (DSHS)

Center for System Coordination and Innovation (CSCI)

P. O. Box 149347, Mail Code 1911, Austin, Texas 78714-9347

Or street address

1100 West 49th Street, Austin, Texas 78756

Fax: Attention: DSHS Public Hearing at (512) 776-7671

**E-mail:** DSHSPublicHearings@dshs.texas.gov

Phone: (512) 776-6206

This forum is open to the public. No reservations are required, and there is no cost to attend. People with disabilities who wish to attend the public hearing and require auxiliary aids or services should call (512) 776-6206 at least 72 hours before the public hearing so appropriate arrangements can be made.

TRD-202401222
Cynthia Hernandez
General Counsel
Department of State Health Services

Filed: March 20, 2024

**\* \* \*** 

Order Placing Zuranolone into Schedule IV, Placing Nine Specific Fentanyl-Related Substances into Schedule I, and Temporarily Placing MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, and 5F-EDMB-PICA, and MMB-FUBICA into Schedule I

The U.S. Drug Enforcement Administration issued an interim final rule placing zuranolone (chemically known as 1-[2-[(3R,5R,8R,9R,10S,13S,14S,17S)-3-hydroxy-3,13-dimethyl-2,4,5,6,7,8,9,10,11,12,14,15,16,17-tetradecahydro-1*H*-cyclopenta[a]phenanthren-17-yl]-2-oxoethyl]pyrazole-4-carbonitrile), including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible, in schedule IV of the Controlled Substances Act effective October 31, 2023. This final rule was published in the *Federal Register*, Volume 88, Number 209, pages 74347-74352.

This scheduling action was taken pursuant to the following:

- 1. Zuranolone has a potential for abuse similar to the drugs or other substances in schedule IV;
- 2. Zuranolone has a currently accepted medical use in treatment in the United States; and,
- 3. Abuse of zuranolone may lead to limited physical dependence or psychological dependence relative to the drugs or other substances in schedule III but similar to other substances in schedule IV.
- The U.S. Drug Enforcement Administration issued a final rule placing the following nine specific fentanyl-related substances, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers whenever the existence of such isomers, esters, and salts is possible, in schedule I of the Controlled Substances Act effective December 7, 2023:
- meta-fluorofentanyl (N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide);
- meta-fluoroisobutyryl fentanyl (N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);
- para-methoxyfuranyl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide);
- 3-furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-phenylfuran-3-carboxamide);
- 2',5'-dimethoxyfentanyl (*N*-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-*N*-phenylpropionamide);
- isovaleryl fentanyl (3-methyl-N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide);

- *ortho*-fluorofuranyl fentanyl (*N*-(2-fluorophenyl)-*N*-(1-phenethylpiperidin-4-yl)furan-2-carboxamide);
- alpha'-methyl butyryl fentanyl (2-methyl-N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide); and
- para-methylcyclopropyl fentanyl (N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)cyclopropanecarboxamide).

This final rule was published in the *Federal Register*, Volume 88, Number 234, pages 85104-85109.

This scheduling action was taken pursuant to the following:

- 1. The abuse potential of *meta*-fluorofentanyl, *meta*-fluoroisobutyryl fentanyl, *para*-methoxyfuranyl fentanyl, 3-furanyl fentanyl, 2',5'-dimethoxyfentanyl, isovaleryl fentanyl, *ortho*-fluorofuranyl fentanyl, *alpha*'-methyl butyryl fentanyl, and *para*-methylcyclopropyl fentanyl is associated with each substance's pharmacological similarity to other schedule I and II mu-opioid receptor agonist substances which have a high potential for abuse. Similar to morphine (schedule II), fentanyl (schedule II), and several schedule I opioid substances that are structurally related to fentanyl, these nine fentanyl-related substances have been shown to bind and act as mu-opioid receptor agonists;
- 2. meta-Fluorofentanyl, meta-fluoroisobutyryl fentanyl, para-methoxyfuranyl fentanyl, 3-furanyl fentanyl, 2',5'-dimethoxyfentanyl, isovaleryl fentanyl, ortho-fluorofuranyl fentanyl, alpha'-methyl butyryl fentanyl, and para-methylcyclopropyl fentanyl, have no currently accepted medical use in treatment in the United States; and
- 3. There is a lack of accepted safety for use of *meta*-fluorofentanyl, *meta*-fluoroisobutyryl fentanyl, para-methoxyfuranyl fentanyl, 3-furanyl fentanyl, 2',5'-dimethoxyfentanyl, isovaleryl fentanyl, *ortho*-fluorofuranyl fentanyl, *alpha'*-methyl butyryl fentanyl, and *para*-methylcyclopropyl fentanyl under medical supervision.
- The U.S. Drug Enforcement Administration issued temporary order placing MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA including their optical and geometric isomer, salts, and salts of isomers, whenever the existence of such isomers and salts is possible, in schedule I of the Controlled Substance Act effective December 12, 2023:

This temporary order was published in the *Federal Register*, Volume 88, Number 237, pages 86040-86046.

This scheduling action was taken pursuant to the following:

- 1. MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA pose an imminent hazard to the public safety;
- 2. MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA have a high potential for abuse;
- 3. MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA have no currently accepted medical use in treatment in the United States; and
- 4. MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA have a lack of accepted safety for use 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, Jennifer Shuford, M.D., does hereby order that the substance zuranolone be placed into schedule IV, nine specific fentanyl-related substances be placed into schedule I, and MDMB-4en-PINACA, 4F-MDMB-BUTICA, ADB-4en-PINACA, CUMYL-PEGACLONE, 5F-EDMB-PICA, and MMB-FUBICA be temporarily placed into schedule I.

#### -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-a-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-a-cetylmethadol, levo-a-acetylmethadol, levomethadyl acetate, or LAAM);
- \*(8) a'-Methyl butyryl fentanyl (2-methyl-N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide);
- (9) a-Methylfentanyl or any other derivative of fentanyl;
- (10)  $\alpha$ -Methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl] N-phenylpropanamide);
- (11) Benzethidine;
- (12)  $\beta$ -Hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide);
- (13)  $\beta$ -Hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide);
- (14) 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);
- (15)  $\beta$ -Methyl fentanyl (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide);
- (16)  $\beta'$ -Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide) (Other name: 3-phenylpropanoyl fentanyl);
- (17) Betaprodine;
- (18) Brorphine (1-(1-(4-bromophenyl)ethyl)piperidin-4-yl)-1,3-dihydro-2H-benzo[d]imidazol-2-one);
- (19) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide);
- (20) Clonitazene:
- (21) Crotonyl fentanyl (Other name: (6-2-5) (E)-N-(1-Phenethylpiperidin-4-yl)-N-phenylbut-2-enamide);
- (22) Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-

Phenylcyclopentanecarboxamide;

- (23) Cyclopropyl fentanyl (*N*-(1-phenethylpiperidin-4-yl)-*N*-phenylcyclopropanecarboxamide);
- (24) Diampromide;
- (25) Diethylthiambutene;
- (26) Difenoxin;
- (27) Dimenoxadol;
- \*(28) 2',5'-Dimethoxyfentanyl (N-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-N-phenylpropionamide);
- (29) Dimethylthiambutene;
- (30) Dioxaphetyl butyrate;
- (31) Dipipanone;
- (32) Ethylmethylthiambutene;
- (33) Etonitazene;

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(34) Etoxeridine;
(35) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-yl)(phenyl)carbamate);
(36) 4-Fluoroisobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)isobutyramide) (Other name: p-fluoroisobutyryl fentanyl);
(37) 2'-Fluoro o-fluorofentanyl (N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-
fluorophenyl)propionamide (Other name: 2'-fluoro 2-fluorofentanyl);
(38) Furanyl fentanyl (N-(1-phenethylpiperdin-4-yl)-N-phenylfuran-2-
carboxamide):
*(39) 3-Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3-
carboxamide);
(40) Furethidine;
(41) Hydroxypethidine:
(42) Isobutyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide;
(43) Isotonitazene (N,N-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1H-
benzimidazol-1-yl)ethan-1-amine);
*(44) Isovaleryl fentanyl (3-methyl-N-(1-phenethylpiperidin-4-yl)-N-
phenylbutanamide);
(45) Ketobemidone;
(46) Levophenacylmorphan;
*(47) m-Fluorofentanyl (N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)propionamide);
*(48) m-Fluoroisobutyryl fentanyl (N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)isobutyramide)
(49) Meprodine;
(50) Methadol;
(51) Methoxyacetyl fentanyl (2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-
phenylacetamide);
(52) 4'-Methyl acetyl fentanyl (N-(1-(4-methylphenethyl)piperidin-4-yl)-N-
phenylacetamide;
(53) 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-
phenylpropanamide);
(54) 3-Methylthiofentanyl (N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-
phenylpropanamide);
(55) Metonitazene (N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-
1-yl)ethan-1-amine);
(56) Moramide;
(57) Morpheridine;
(58) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
(59) MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
(60) Noracymethadol;
(61) Norlevorphanol;
(62) Normethadone;
(63) Norpipanone;
(64) Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-
yl)acetamide);
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(65) o-Fluoroacryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
vl)acrylamide);
(66) o-Fluorobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)butyramide (Other name:2-fluorobutyryl fentanyl);
(67) o-Fluorofentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)propionamide) (Other name: 2-fluorofentanyl);
*(68) o-Fluorofuranyl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)furan-2-carboxamide);
(69) o-Fluoroisobutyryl fentanyl (N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)isobutyramide);
(70) o-Methyl acetylfentanyl (N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-
yl)acetamide (Other name: 2-methyl acetylfentanyl);
(71) o-Methyl methoxyacetyl fentanyl (2-methoxy-N-(2-methylphenyl)-N-(1-
phenethylpiperidin-4-yl)acetamide (Other name: 2-methyl methoxyacetyl
fentanyl);
(72) p-Chloroisobutyryl fentanyl (N-(4-chlorophenyl)-N-(1- phenethylpiperidin-4-
yl)isobutyramide;
(73) p-Fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)butyramide);
(74) p-Fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4 piperidinyl]
propanamide);
(75) p-Fluoro furanyl fentanyl (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-
yl)furan-2-carboxamide);
(76) p-Methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1- phenethylpiperidin-4-
yl)butyramide;
*(77) p-Methoxyfuranyl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-
vI)furan-2-carboxamide);
*(78) p-Methylcyclopropyl fentanyl (N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-
yl)cyclopropanecarboxamide);
(79) p-Methylfentanyl (N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-
yl)propionamide (Other name: 4-methylfentanyl);
(80) PEPAP (1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
(81) Phenadoxone;
(82) Phenampromide;
(83) Phencyclidine;
(84) Phenomorphan;
(85) Phenoperidine;
(86) Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide (Other
name: benzoyl fentanyl);
(87) Piritramide;
(88) Proheptazine;
(89) Properidine;
(90) Propiram;
(91) Tetrahydrofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-
phenyltetrahydrofuran-2-carboxamide);
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- (92) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);
- (93) Thiofuranyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide (Other names: 2-thiofuranyl fentanyl; thiophene fentanyl); (94) Tilidine;
- (95) Trimeperidine;
- (96) U-47700 (3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide;
- (97) Valeryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide); and,
- (98) Zipeprol (1-methoxy-3-[4-(2-methoxy-2-phenylethyl)piperazin-1-yl]-1-phenylpropan-2-ol).

# -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, alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups,

(1-1-3) Substitution in or on the piperidine ring with alkyl,

alkenyl, alkoxyl, 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:  $\beta'$ -phenyl fentanyl; hydrocinnamoyl fentanyl); and,

- (1-2-4) *N*-(1-Phenethylpiperidin-4-yl)-*N*-phenylthiophene-2-carboxamide (Other name: thiofuranyl fentanyl).
- (2) 2-(2-(4-Butoxybenzyl)-5-nitro-1*H*-benzimidazol-1-yl)-*N*,*N*-diethylethan-1-amine (Other name: butonitazene);
- (3) 2-(2-(4-Ethoxybenzyl)-1*H*-benzimidazol-1-yl)-*N*,*N*-diethylethan-1-amine (Other names: etodesnitazene; etazene);
- (4) N,N-Diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1- yl)ethan-1-amine (Other name: flunitazene);
- (5) N,N-Diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine (Other name: metodesnitazene);
- (6) 2-(4-Ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole (Other names: *N*-pyrrolidino etonitazene; etonitazepyne);
- (7) *N*,*N*-Diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1*H*-benzimidazol-1-yl)ethan-1-amine (Other name: protonitazene);
- (8) 4-(2-chlorophenyl)-2-ethyl-9-methyl-6H-thieno[3,2-f][1,2,4]triazolo[4,3-a][1,4]diazepine (Other name: etizolam);
- (9) 8-chloro-6-(2-fluorophenyl)-1-methyl-4H-benzo[f][1,2,4]triazolo[4,3-a][1,4]diazepine (Other name: flualprazolam);
- (10) 6-(2-chlorophenyl)-1-methyl-8-nitro-4H-benzo[f][1,2,4]triazolo[4,3-a][1,4]diazepine (Other name: clonazolam);
- (11) 8-bromo-6-(2-fluorophenyl)-1-methyl-4H-benzo[f][1,2,4]triazolo[4,3- $\alpha$ ][1,4]diazepine (Other names: 8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3- $\alpha$ ][1,4]benzodiazepine and flubromazolam);
  - (12) 7-chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2*H*-
- benzo[e][1,4]diazepin-2-one (Other name: diclazepam);
- \*(13) Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1*H*-indazole-3-carboxamido)butanoate (Other name: MDMB-4en-PINACA)
- \*(14) Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate (Other names: 4F-MDMB-BUTICA; 4F-MDMB-BICA)
- \*(15) N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-en-1-yl)-1H-indazole-3-carboxamide (Other name: ADB-4en-PINACA)
- \*(16) 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido[4,3-b]indol-1-one (Other names: CUMYL-PEGACLONE; SGT-151)
- \*(17) Ethyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-dimethylbutanoate (Other names: 5F-EDMB-PICA; 5F-EDMB-2201)
- \*(18) Methyl 2-(1-(4-fluorobenzyl)-1*H*-indole-3-carboxamido)-3-methyl butanoate (Other name: MMB-FUBICA)

## -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:

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Alfaxalone (5a-pregnan-3a-ol-11,20-dione);
      (2) Alprazolam;
      (3) Barbital;
      (4) Brexanolone (3a-hydroxy-5a-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 (Other name: mephobarbital);
      (39) Midazolam;
      (40) Nimetazepam;
      (41) Nitrazepam;
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- (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;
- (58) Zopiclone; and
- \*(59) Zuranolone.

# Changes are marked by an asterisk(\*)

TRD-202401216 Cynthia Hernandez General Counsel

Department of State Health Services

Filed: March 19, 2024

### Texas Department of Housing and Community Affairs

Notice of Public Hearings and Public Comment Period on the Draft 2024 State of Texas Analysis of Impediments to Fair Housing Choice

The Texas Department of Housing and Community Affairs (TDHCA) will hold four public hearings during a 31-day Public Comment period to accept public comment on the draft 2024 State of Texas Analysis of Impediments to Fair Housing Choice (AI). The Public Comment period for the draft AI will be held Wednesday, April 3, 2024 - Friday, May 3, 2024.

The public hearings for the 2024 AI are scheduled to take place as follows:

Tuesday, April 16, 2024 - 6:30 p.m.

Carver Branch, Austin Public Library

Room #1

1161 Angelina Street

Austin, Texas 78702

Thursday, April 18, 2024 - 6:00 p.m.

Via GoToWebinar

https://attendee.gotowebinar.com/register/1281746222460736607

Dial-in number: +1 (562) 247-8422, Access Code 583-444-377 (persons who use the dial-in number and access code without registering online will only be able to hear the public hearing and will not be able to ask questions or provide comments. Staff recommend that attendees without computer access use the GoToWebinar phone app in order to participate fully.)

Tuesday, April 23, 2024 - 12:30 p.m.

come. dream. come. build.

Father Armand Matthew Meeting Room

901 E Levee Street

Brownsville, Texas 78520

Thursday, April 25, 2024 - 11:00 a.m.

J. Erik Jonsson Central Library

1st Floor Auditorium

1515 Young Street

Dallas, Texas 75201

The State of Texas is releasing for public comment the 2024 Analysis of Impediments to Fair Housing Choice (AI) planning document, which guides the state's affordable housing strategy for the next five years. This is a process required of recipients of U.S. Department of Housing and Urban Development (HUD) Community Planning and Development grant funds as part of their obligation to undertake affirmatively further fair housing (AFFH) under the Fair Housing Act. The Texas Department of Housing and Community Affairs (TDHCA) is the agency in Texas that is primarily responsible for creating the AI.