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# MISSOUR, המומומומומומומומ ΠП SALUS POPULI SUPREMA LEX ESTO "The welfare of the people shall be the supreme law" пппппппп ПП Π Π пп REGISTER

John R. Ashcroft 🛞 Secretary of State

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# MISSOURI

Department of Health and Senior Services



# REGISTER

October 17, 2022

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June 1, 2022	July 1, 2022	July 31, 2022	August 30, 2022
June 15, 2022	July 15, 2022	July 31, 2022	August 30, 2022
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January 3, 2023	February 1, 2023	February 28, 2023	March 30, 2023
January 17, 2023	February 15, 2023	February 28, 2023	March 30, 2023

Documents will be accepted for filing on all regular workdays from 8:00 a.m. until 5:00 p.m. We encourage early filings to facilitate the timely publication of the *Missouri Register*. Orders of Rulemaking appearing in the *Missouri Register* will be published in the *Code of State Regulations* and become effective as listed in the chart above. Advance notice of large volume filings will facilitate their timely publication. We reserve the right to change the schedule due to special circumstances. Please check the latest publication to verify that no changes have been made in this schedule. To review the entire year's schedule, please see the website at sos.mo.gov/adrules/pubsched.

## HOW TO CITE RULES AND RSMO

#### RULES

The rules are codified in the Code of State Regulations in this system-

Title	CSR	Division	Chapter	Rule
3	Code of	10-	4	115
Department	State	Agency	General area	Specific area
	Regulations	division	regulated	regulated

and should be cited in this manner: 3 CSR 10-4.115.

Each department of state government is assigned a title. Each agency or division in the department is assigned a division number. The agency then groups its rules into general subject matter areas called chapters and specific areas called rules. Within a rule, the first breakdown is called a section and is designated as (1). Subsection is (A) with further breakdown into paragraphs 1., subparagraphs A., parts (I), subparts (a), items I. and subitems a.

The rule is properly cited by using the full citation; for example, 3 CSR 10-4.115, NOT Rule 10-4.115.

Citations of RSMo are to the Missouri Revised Statutes as of the date indicated.

### Code and Register on the Internet

The Code of State Regulations and Missouri Register are available on the Internet.

The Code address is sos.mo.gov/adrules/csr/csr

The *Register* address is <u>sos.mo.gov/adrules/moreg/moreg</u>

These websites contain rulemakings and regulations as they appear in the Code and Registers.

#### MISSOURI REGISTER

Rules appearing under this heading are filed under the authority granted by section 536.025, RSMo. An emergency rule may be adopted by an agency if the agency finds that an immediate danger to the public health, safety, or welfare, or a compelling governmental interest requires emergency action; follows procedures best calculated to assure fairness to all interested persons and parties under the circumstances; follows procedures which comply with the protections extended by the Missouri and the United States Constitutions; limits the scope of such rule to the circumstances creating an emergency and requiring emergency procedure, and at the time of or prior to the adoption of such rule files with the secretary of state the text of the rule together with the specific facts, reasons, and findings which support its conclusion that there is an immediate danger to the public health, safety, or welfare which can be met only through the adoption of such rule and its reasons for concluding that the procedure employed is fair to all interested persons and parties under the circumstances.

Rules filed as emergency rules may be effective not less than ten (10) business days after filing or at such later date as may be specified in the rule and may be terminated at any time by the state agency by filing an order with the secretary of state fixing the date of such termination, which order shall be published by the secretary of state in the Missouri Register as soon as practicable.

All emergency rules must state the period during which they are in effect, and in no case can they be in effect more than one hundred eighty (180) calendar days or thirty (30) legislative days, whichever period is longer. Emergency rules are not renewable, although an agency may at any time adopt an identical rule under the normal rulemaking procedures.

#### Title 19 – DEPARTMENT OF HEALTH AND SENIOR SERVICES Division 30 – Division of Regulation and Licensure Chapter 1 – Controlled Substances

#### **EMERGENCY AMENDMENT**

**19 CSR 30-1.002 Schedules of Controlled Substances.** The department is amending section (1).

PURPOSE: This amendment updates the Schedules of Controlled Substances to be consistent with 21 CFR Part 1308.

EMERGENCY STATEMENT: The United States Department of Justice Drug Enforcement Administration (DEA) continually evaluates substances to determine their clinical application and potential for abuse. Based on their evaluation, the DEA issues scheduling actions to place substances in the appropriate controlled substance schedules. The majority of these scheduling actions consist of temporarily and permanently scheduling newly-discovered illicit substances in Schedule I. Proper scheduling of these substances allow law enforcement to take action to prevent the further distribution of these substances. Scheduling substances in Schedules II-V allows practitioners to be informed about the potential for addiction/abuse of the substances and prescribe the substances appropriately. Section 195.015, RSMo charges the department with similarly controlling substances as they are controlled under federal law. Section

195.015.4 requires the Department of Health and Senior Services to submit emergency rules to the Secretary of State within thirty days of a federal scheduling action to allow for similar inclusion, rescheduling, or deletion of controlled substances with this schedule. While this time frame is difficult to achieve given the various approvals and reviews needed prior to the Department scheduling any rule with the Secretary of State, the Department still acts to effectuate these scheduling actions as quickly as possible. This emergency amendment includes all federal scheduling actions since the last amendment of this rule in 2020. This emergency amendment is necessary to protect Missouri's governmental interest in keeping its controlled substances schedules up-to-date as much as practically possible in order to protect its citizens and to aid law enforcement in its prosecution of those who illegally distribute these substances. As a result, the Department of Health and Senior Services finds a compelling governmental interest, which requires this emergency action. A proposed amendment, which covers the same material, is published in this issue of the Missouri Register. The scope of this emergency amendment is limited to the circumstances creating the emergency and complies with the protections extended in the Missouri and United States Constitutions. The Department of Health and Senior Services believes this emergency amendment is fair to all interested persons and parties under the circumstances. Subject to section 536.025, this emergency amendment was filed September 12, 2022, becomes effective October 3, 2022, and expires March 31, 2023.

(1) Schedules of Controlled Substances.

(A) Schedule I shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Drug Enforcement Administration (DEA) Controlled Substances Code Number set forth opposite it.

1. Opiates. Unless specifically excepted or unless listed in another schedule, any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation:

A. Acetyl-alpha-methylfentanyl	
(N-(1-(1-methyl-2-phenethyl)-	
4-piperidinyl)-N-	
phenylacetamide)	9815
B. Acetylmethadol	9601
C. Acetyl fentanyl (N-(1-	
phenethylpiperidin-4-yl)-	
N-phenylacetamide)	9821
D. N-(1-phenethylpiperidin-	
4-yl)-N-phenylacrylamide,	
its isomers, esters, ethers,	
salts, and salts of isomers,	
esters, and ethers (Other	
names: acryl fentanyl,	
acryloylfentanyl)	9811
E. AH-7921(3,4-dichloro-	
N-[(1-dimethylamino)	
cyclohexylmethyl]	
benzamide)	9551
F. Allylprodine	9602
G. Alphacetylmethadol (except	
levoalphacetylmethadol	
also known as levo-alpha-	
acetylmethadol levothadyl	
acetate or LAAM)	9603
H. Alphameprodine	9604

# **EMERGENCY RULES**

I. Alphamethadol	9605	LL. Fentanyl carbamate (ethyl	
J. Alpha-methylfentanyl		(1-phenethylpiperidin-4-yl)	
(N-1-(alphamethyl-beta-		(phenyl)carbamate)	9851
phenyl) ethyl-4-piperidyl)		MM. N-(4-fluorophenyl)-N-	
propionanilide; 1-(1-methyl-		(1-phenethylpiperidin-4-	
2-phenylethyl)-4 ((N-		yl)isobutyramide, its	
propanilido) piperidine)	9814	isomers, esters, ethers,	
K. Alpha-methylthiofentanyl		salts, and salts of isomers,	
(N-(1-methyl-2-(2-thienyl)		esters, and ethers (Other	
ethyl-4-piperidinyl)-N-		names: 4-fluoroisobutyryl	
phenylpropanamide)	9832	fentanyl, para-	
L. Benzethidine	9606	fluoroisobutyryl	
M. Betacetylmethadol	9607	fentanyl)	9824
N. Beta-hydroxyfentanyl		NN. 2'-Fluoro ortho-	
(N-(1-(2-hydroxy-2-		fluorofentanyl (N-(1-(2-	
phenethyl)-4-piperidinyl)-		fluorophenethyl) piperidin-	
N-phenylpropanamide)	9830	4-yl)-N-(2-fluorophenyl)	
O. Beta-hydroxy-3-		propionamide (Other names:	
methylfentanyl (other name:		2'-fluoro 2-	
N-(1-(2-hydroxy-2-phenethyl)-		fluorofentanyl)	9855
3-methyl-4-piperidinyl)-N-		OO. N-(1-phenethylpiperidin-	
phenylpropanamide)	9831	4-yl)-N-phenylfuran-2-	
P. N-[1-[2-hydroxy-2-(thiophen-		carboxamide (Other names:	
2-yl) ethyl]piperidin-4-yl]-		furanyl fentanyl)	9834
N-phenylpropionamide		PP. Furethidine	9626
(Other names:		QQ. Hydroxypethidine	9627
beta-hydroxythiofentanyl)	9836	RR. N-(1-phenethylpiperidin-	
Q. Betameprodine	9608	4-yl)-N-phenylisobutyramide	
R. Betamethadol	9609	(Other name: isobutyryl fentanyl)	9827
S. beta-Methyl fentanyl		SS. Isotonitazene ( <i>N</i> , <i>N</i> -diethyl-2-(2-	
(N-phenyl-N-(1-(2-		(4-isopropoxybenzyl)-5-nitro-	
phenylpropyl)piperidin-4-yl)		1 <i>H</i> -benzimidazol-1-yl)	
propionamide (Other name:		ethan-1-amine)	9614
β-methyl fentanyl)	9856	[SS.]TT. Ketobemidone	9628
T. beta'-Phenyl fentanyl		[TT.]UU. Levomoramide	9629
(N-(1-phenethylpiperidin-4-yl)-		[UU.]VV. Levophenacylmorphan	9631
N,3-diphenylpropanamide		[VV.]WW. Methoxyacetyl fentanyl	
(Other names: $\beta'$ -phenyl		(2-methoxy-N-(1-	
fentanyl; 3-phenylpropanoyl		phenethylpiperidin-4-yl)-	
fentanyl)	9842	N-phenylacetamide	9825
U. Betaprodine	9611	[WW.]XX. 4'-Methyl acetyl fentanyl	
V. Clonitazene	9612	(N-(1-(4-methylphenethyl)	
W. Crotonyl fentanyl ((E)-N-(1-		piperidin-4-yl)-N-	
phenethylpiperidin-4-yl)-N-		phenylacetamide)	9819
phenylbut-2-enamide)	9844	[XX.]YY. 3-Methylfentanyl (N-(3-	
X. N-(1-phenethylpiperidin-		methyl-1-(2-phenylethyl)-4-	
4-yl)-N-		piperidyl)-N-	
Phenylcyclopentanecarboxamide		phenylproanamide), its	
(Other name: cyclopentyl		optical and geometric	
fentanyl)	9847	isomers, salts, and salts	
Y. Cyclopropyl fentanyl (N-(1-	001,	of isomers	9813
phenethylpiperidin-4-yl)-N-		[YY.]ZZ. 3-Methylthiofentanyl (N-	0010
phenylcyclopropanecar-		(3-methyl-1-(2-	
boxamide)	9845	thienyl)ethyl-4-piperidinyl)-	
Z. Dextromoramide	9613	N-phenylpropanamide)	9833
AA. Diampromide	9615	[ZZ.]AAA. Morpheridine	9632
BB. Diethylthiambutene	9616	[AAA.]BBB. MPPP (1-methyl-4-	3001
CC. Difenoxin	9168	phenyl-4-	
DD. Dimenoxadol	9617	propionoxypiperidine)	9661
EE. Dimepheptanol	9618	[BBB.]CCC. MT-45 (1-cyclohexyl-	5001
FF. Dimethylthiambutene	9619	4-(1,2-diphenylethyl)	
GG. Dioxaphetyl butyrate	9621	piperazine)	(9560)
HH. Dipipanone	9622	[CCC.]DDD. Noracymethadol	9633
II. Ethylmethylthiambutene	9623	[DDD.]EEE. Norlevorphanol	9634
II. Etonitazene	9624	[EEE.]FFF. Normethadone	9635
JJ. Etomtazene KK. Etoxeridine	9625	[FFF.]GGG. Norpipanone	9636
	5020	L' I Jooo. Norpipullone	5050

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[GGG.]HHH. N-(2-fluorophenyl)-2- methoxy-N-(1-		[SSS.]TTT. para-Methylfentanyl (N-(4-methylphenyl)-N-	
phenethylpiperidin-4-		(1-phenethylpiperidin-4-yl)	
yl)acetamide, its isomers,		propionamide (Other	
esters, ethers, salts, and		Name: 4-	
salts of isomers, esters,		methylfentanyl)	9817
and ethers (Other name:		[TTT.]UUU. PEPAP (1-(-2-phenethyl)-	
ocfentanil)	9838	4-phenyl-4-	
[HHH.]III. ortho-Fluoroacryl		acetoxypiperidine)	9663
fentanyl (N-(2-		[UUU.]VVV. Phenadoxone	9637
fluorophenyl)-N-(1-		[VVV.]WWW. Phenampromide	9638
phenethylpiperidin-4-yl)		[WWW.]XXX. Phenomorphan	9647 9641
acrylamide)	9852	[XXX.]YYY. Phenoperidine [YYY.]ZZZ. Phenyl fentanyl (N-(1-	9041
[///.]]]]. ortho-Fluorobutyryl fentanyl		phenethylpiperidin-4-yl)-	
(N-(2-fluorophenyl)-N-(1-		N-phenylbenzamide	
phenethylpiperidin-4-yl) buturamida (Othor Name:		(Other name: benzoyl	
butyramide (Other Name: 2-fluorobutyryl fentanyl)	9846	fentanyl)	9841
[JJJ.]KKK. ortho-Fluorofentanyl (N-(2-	9640	[ZZZ.]AAAA. Piritramide	9642
fluorophenyl)-N-(1-		[AAAA.]BBBB. Proheptazine	9643
phenethylpiperidin-4-yl)		[BBBB.]CCCC. Properidine	9644
propionamide); other name:		[CCCC.]DDDD. Propiram	9649
2-fluorofentanyl)	9816	[DDDD]EEEE. Racemoramide	9645
[KKK.]LLL. ortho-Fluoroisobutyryl	5010	<i>[EEEE.]</i> FFFF. N-(1-	
fentanyl (N-(2-		phenethylpiperidin-4-yl)-	
fluorophenyl)-N-(1-		N-phenyltetrahydrofuran-	
phenethylpiperidin-4-		2-carboxamide, its	
yl)isobutyramide)	9853	isomers, esters, ethers,	
[LLL.]MMM. ortho-Methyl acetylfentanyl		salts, and salts of isomers,	
(N-(2-methylphenyl)-N-(1-		esters, and ethers (Other name: tetrahydrofuranyl	
phenethylpiperidin-4-yl)		fentanyl)	9843
acetamide (Other name:		[FFFF.]GGGG. Thiofentany (N-phenyl-	3043
2-methyl acetylfentanyl)	9848	N-(1-(2-thienyl)ethyl-4-	
[MMM.]NNN. ortho-Methyl		piperidinyl)-	
methoxyacetyl fentanyl		propanamide	9835
(2-methoxy-N-(2-		[GGGG.]HHHH. Thiofuranyl fentanyl	
methylphenyl)-N-(1-		(N-(1-phenethylpiperidin-	
phenethylpiperidin-4-yl)		4-yl)-N-phenylthiophene-	
acetamide (Other name: 2-methyl		2-carboxamide (Other	
methoxyacetyl		names: 2-thiofuranyl	
fentanyl)	9820	fentanyl; thiophene	
[NNN.]OOO. N-(4-chlorophenyl)-N-	3020	fentanyl)	9839
(1-phenethylpiperidin-		[HHHH.]IIII. Tilidine [////.]IIII. Trimeperidine	9750
4-yl)isobutyramide		2. Opium derivatives. Unless specifically ex	9646 conted or
(Other name: para-		unless listed in another schedule, any of the follow	
chloroisobutyryl		derivatives, its salts, isomers, and salts of isomers	
fentanyl)	9826	the existence of such salts, isomers, and salts of	
[OOO.]PPP. para-Fluorobutyryl		possible within the specific chemical designation:	
fentanyl (N-(4-		A. Acetorphine	9319
fluorophenyl)-N-(1-		B. Acetyldihydrocodeine	9051
phenethylpiperidin-4-		C. Benzylmorphine	9052
yl)butyramide)	9823	D. Codeine methylbromide	9070
[PPP.]QQQ. Para-fluorofentanyl(N-		E. Codeine-N-Oxide	9053
(4-fluorophenyl)-N-(1-(2- phenethyl)-4-piperidinyl)		F. Cyprenorphine	9054
propanamide	9812	G. Desomorphine	9055
[QQQ.]RRR. para-Fluoro furanyl	9012	H. Dihydromorphine I. Drotebanol	9145 9335
fentanyl (N-(4-		J. Etorphine (except	9333
fluorophenyl)-N-(1-		hydrochloride salt)	9056
phenethylpiperidin-4-		K. Heroin	9030
yl)furan-2-carboxamide)	9854	L. Hydromorphinol	9301
[RRR.]SSS. para-Methoxybutyryl		M. Methyldesorphine	9302
fentanyl (N-(4-		N. Methyldihydromorphine	9304
methoxyphenyl)-N-(1-		O. Morphine methylbromide	9305
phenethylpiperidin-4-yl)		P. Morphine methylsulfonate	9306
butyramide)	9837	Q. Morphine-N-Oxide	9307

# **EMERGENCY RULES**

7249

	R. Myrop	hine			9308
	S. Nicoco				9309
	T. Nicom	orphine			9312
	U. Normo	orphine			9313
	V. Pholco	dine			9314
	W. Theba	con			9315
З	Oniato	Similar	Synthetic	Substances	Substances

3. Opiate Similar Synthetic Substances. Substances scheduled by the United States Drug Enforcement Administration as substances that share a pharmacological profile similar to fentanyl, morphine, and other synthetic opioids, unless specifically excepted or unless listed in another schedule. These substances are –

A. Butyryl fentanyl (N-	
(1-phenethylpiperidin-4-yl)-	
N-phenylbutyramide)	9822
B. U–47700 (3,4-Dichloro-	
N-[2-(dimethylamino)	
cyclohexyl]-N-	
methylbenzamide)	9547
C. N-(1-phenethylpiperidin-	
4-yl)-N-phenylpentanamide	
(Other name: valeryl	
fentanyl)	9840

4. Hallucinogenic substances. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or preparation, which contains any quantity of the following hallucinogenic substances or which contains any of its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation (For purposes of paragraph (1)(A)4. of this rule only, the term isomer includes the optical, position, and geometric isomers.):

A. Alpha-ethyltryptamine

Some trade or other names: etryptamine; Monase; alpha-ethyl-1H-indole-3-ethenamine; 3-(2-aminobutyl)indole; alpha-ET; and AET;

B. 4-bromo-2,5-	
dimethoxyamphetamine	7391
Some trade or other names: 4-bromo-2, 5- di	methoxy-a-
methylphenethylamine; 4-bromo- 2, 5-DMA;	
C 4 brome 2 F	

C. 4-bromo-2,5-		
dimethoxyphenethylamine		7392
D. 2,5-dimethoxyamphetamir	ie	7396
Some trade or other	names:	2,5-dimethoxy-
amethylphenethylamine; 2,5-DMA;		
E. 2,5-dimethoxy-4-		
ethylamphetamine		7399
Some trade or other names: DOET;		
F. 2,5-dimethoxy-4-(n)-		
propylthiophenethylamine		
(other name: 2C-T-7)		7348
G. 2-(2,5-Dimethoxy-4-(n)-		
propylphenyl) ethanamine		
(2C-P)		7524
H. 2-(2,5-Dimethoxy-4-		
ethylphenyl) ethanamine		
(2C-E)		7509
I. 2-(2,5-Dimethoxy-4-		
methylphenyl) ethanamine		
(2C-D)		7508
J. 2-(2,5-Dimethoxy-4-nitro-		
phenyl) ethanamine		
(2C-N)		7521
K. 2-(2,5-Dimethoxyphenyl)		
ethanamine (2C-H)		7517
L. 2-(4-Chloro-2,5-		
dimethoxyphenyl)		
ethanamine (2C-C)		7519

	,
M. 2-(4-Ethylthio-2,5-	
dimethoxyphenyl)	
ethanamine (2C-T-2)	7385
N. 2-(4-Iodo-2,5-	
dimethoxyphenyl)	
ethanamine (2C-I)	7518
O. 2-(4-Isopropylthio)-2,5-	
dimethoxyphenyl)	
ethanamine (2C-T-4)	7532
P. 4-methoxyamphetamine	7411
Some trade or other names: 4-methoxy-	
amethylphenethylamine; paramethoxyamphetamin	ie; PMA;
Q. 5-methoxy-3,4-	
methylenedioxyamphetamine	7401
R. 4-methyl-2,5-	7005
dimethoxyamphetamine	7395
Some trade and other names: 4-methyl-2, 5- din	netnoxy-a-
methylphenethylamine; DOM; and STP;	
S. 3,4-	7400
methylenedioxyamphetamine	7400
T. 3,4-methylenedioxymetham- phetamine(MDMA)	7405
U. 3,4-methylenedioxy-N-	7405
ethylamphetamine (also	
known as N-ethylalpha-	
methyl-3,4 (methylenedioxy)	
phenethylamine, N-ethyl	
MDA, MDE, and MDEA)	7404
V. N-hydroxy-3,4-	, 10 1
methylenedioxyamphetamine	
(also known as N-hydroxy-	
alpha-methyl-3,4	
(methylenedioxy)	
phenethylamine and N-	
hydroxy MDA)	7402
W. 3,4,5-	
trimethoxyamphetamine	7390
X. 5-MeO-DMT or 5-methoxy-	
N,N-dimethyltryptamine	7431
Y. Alpha-methyltryptamine	7432
Z. Bufotenine	7433
Some trade and other names: 3-(b-Dimethylam	ninoethyl)-
5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-ind	
N-dimethylserotonin; 5-hydroxy-N, N-dimethyltr	yptamine;
mappine;	
AA. Diethyltryptamine	7434
Some trade and other names: N, N-Diethyltr	yptamine;
DET;	- 40-
BB. Dimethyltryptamine	7435
Some trade or other names: DMT;	
CC. 5-methoxy-N,N-	
diisopropyltryptamine	
(other name: 5-	7420
(other name: 5- MeODIPT)	7439
(other name: 5- MeODIPT) DD. Ibogaine	7260
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8	7260 ,9,10,12,13-
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido	7260 ,9,10,12,13-
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga;	7260 ,9,10,12,13- [1',2':1,2]
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide	7260 ,9,10,12,13- [1',2':1,2] 7315
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana	7260 ,9,10,12,13- [1',2':1,2]
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana;	7260 ,9,10,12,13- [1',2':1,2] 7315 7360
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline	7260 ,9,10,12,13- [1',2':1,2] 7315 7360 7381
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline HH. Parahexyl	7260 9,10,12,13- [1',2':1,2] 7315 7360 7381 7374
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline HH. Parahexyl Some trade or other names: 3-Hexyl-1- hydrox	7260 9,9,10,12,13- [1',2':1,2] 7315 7360 7381 7374 xy-7,8,9,10-
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline HH. Parahexyl Some trade or other names: 3-Hexyl-1- hydrox tetrahydro-6,6,9-trimethyl- 6H-dibenzo[b,d]pyran; Sy	7260 9,9,10,12,13- [1',2':1,2] 7315 7360 7381 7374 xy-7,8,9,10-
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline HH. Parahexyl Some trade or other names: 3-Hexyl-1- hydrox tetrahydro-6,6,9-trimethyl- 6H-dibenzo[b,d]pyran; Sy II. Peyote	7260 ,9,10,12,13- [1',2':1,2] 7315 7360 7381 7374 xy-7,8,9,10- mhexyl; 7415
(other name: 5- MeODIPT) DD. Ibogaine Some trade and other names: 7-Ethyl- 6,6ß,7,8 octahydro-2-methoxy-6, 9-methano-5H-pyrido azepino[5,4-b] indole; Tabernanthe iboga; EE. Lysergic acid diethylamide FF. Marihuana Some trade or other names: marijuana; GG. Mescaline HH. Parahexyl Some trade or other names: 3-Hexyl-1- hydrox tetrahydro-6,6,9-trimethyl- 6H-dibenzo[b,d]pyran; Sy	7260 9,9,10,12,13- [1',2':1,2] 7315 7360 7381 7374 xy-7,8,9,10- mhexyl; 7415 potanically

the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salt, derivative, mixture, or preparation of such plant, its seeds, or extracts;

	-,
JJ. N-ethyl-3-piperidyl	
benzilate	7482
KK. N-methyl-3-piperidyl	
benzilate	7484
LL. Psilocybin	7437
MM. Psilocyn	7438
NN Tetrohudroconnohinolo noturollu	contained in a

NN. Tetrahydrocannabinols naturally contained in a plant of the genus Cannabis (cannabis 7370 plant), 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, or both, with similar chemical structure and pharmacological activity to those substances contained in the plant, such as the following:

(I) 1 cis or trans tetrahydrocannabinol and their optical isomers;

(II) 6 cis or trans tetrahydrocannabinol and their optical isomers;

(III) 3,4 cis or trans tetrahydrocannabinol and its optical isomers; and

(IV) Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions are covered;

covered,	
OO. Ethylamine analog of	
phencyclidine	7455
Some trade or other names: N-ethyl-1- phenylcyclohexyla	
(1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclo	hexyl)-
ethylamine, cyclohexamine, PCE;	
PP. Pyrrolidine analog of	
phencyclidine	7458
Some trade or other names: 1-(1-phenylcyclohexyl)-pyrro	olidine
PCPy, PHP;	
QQ. Thiophene analog of	
phencyclidine	7470
Some trade or other names: 1-(1-(2-thienyl)- cyclo	hexvl)-
piperidine, 2-thienyl analog of phencyclidine, TPCP, TCF	
RR. 1-(1-(2-thienyl)cyclohexyl)	
pyrrolidine	7473
Some other names: TCPy;	
SS. Salvia divinorum	
TT. Salvinorin A	
UU. 3-Fluoromethcathinone	1233
VV. 4-Fluoromethcathinone	1238
WW. Mephedrone, or 4-	1200
methylmethcathinone	1248
XX. Methylenedioxy-	12 10
pyrovalerone, MDPV, or	
(1-(1,3-Benzodioxol-5-yl)-	
2-(1-pyrrolidinyl)-1-	
pentanone	7535
YY. Methylone, or 3,4-	1000
Methylenedioxy-	
methcathinone	7540
ZZ. Quinolin-8-yl 1-pentyl-	7540
1Hindole-3-carboxylate	
(PB-22; QUPIC)	7222
AAA. Quinolin-8-yl 1-(5-	1222
fluoropentyl)-1H-indole-	
3-carboxylate (5-fluoro-	
PB-22; 5F-PB-22)	7225
PD-22, JF-PD-22)	7223
BBB. N-(1-amino-3-methyl-1-	
oxobutan-2-yl)-1- (4 fluorobongyl) I Uindagolo	
(4-fluorobenzyl)-1Hindazole-	
3-carboxamide (AB-	7010
FUBINACA)	7012

CCC. N-(1-amino-3, 3-dimethyl-1-	
oxobutan-2-yl)-1-pentyl-	
1H-indazole-3-carboxamide	
(ADB-PINACA)	7035
DDD. (1-pentyl-1H-indol-3-yl)	
(2,2,3,3-	
tetramethylcyclopropyl)	
methanone (Other names:	
UR-144, 1-pentyl-3-	
(2,2,3,3-	
tetramethylcyclopro-	
poyl)indole)	7144
EEE. [1-(5-fluoro-pentyl)-	
1Hindol-3-yl](2,2,3,3-	
tetramethylcyclopropyl)	
methanone (Other names:	
5-fluoro-UR-144, 5-F-UR-	
144, XLR11, 1-(5-fluo-	
ropentyl)-3-(2,2,3,3-	
tetramethylcyclopro-	
poyl)indole)	7011
FFF. N-(1-adamantyl)-1-pentyl-	
1Hindazole-3-carboxamide	
(Other names: APINACA,	
AKB48)	7048
GGG. 2-(4-iodo-2,5-	
dimethoxyphenyl)-N-(2-	
methoxybenzyl)ethanamine	
(Other names: 251-	
NBOMe; 2C-I-NBOMe;	
25I; Cimbi-5)	7538
HHH. 2-(4-chloro-2,5-	
dimethoxyphenyl)-N-(2-	
methoxybenzyl)ethanamine	
(Other names: 25C-	
NBOMe; 2C-C-NBOMe;	7507
25C; Cimbi-82)	7537
III. 2-(4-bromo-2,5- dimethoxyphenyl)-N-(2-	
methoxybenzyl)ethanamine	
(Other names: 25B-	
NBOMe; 2C-B-NBOMe;	
25B; Cimbi-36)	7536
III. 4-methyl-N-ethylcathinone	7550
(Other names: 4-MEC; 2-	
(ethylamino)-1-(4-	
methylphenyl)propan-	
1-one)	1249
KKK. 4-methyl-alphapyrrolid-	12 10
inopropiophenone,	
(Other names: 4-MePPP;	
MePPP; 4-methyl-	
α-pyrrolidinopropiophenone;	
1-(4-methylphenyl)-2-	
(pyrrolidin-1-yl)-propan-	
1-one)	7498
LLL. alphapyrrolidinopentio-	
phenone	
Other names: α-PVP; α-	
pyrrolidinovalerophenone;	
1-phenyl-2-(pyrrolidin-1-	
yl)pentan-1-one)	7545
MMM. Butylone	
(Other names: bk-	
MBDB; 1-(1,3-	
benzodioxol-5-	
yl)-2-	
(methylamino)butan-1-	
one)	7541

# **EMERGENCY RULES**

NNN. Pentedrone		YYY. N-(1-amino-3,3-dimethyl-	
(Other names: α-		1-oxobutan-2-yl)-1-(4-	
methylaminovalerophenone;		fluorobenzyl)-1H-indazole-	
2-(methylamino)-1-		3-carboxamide	
phenylpentan-1-one)	1246	(Other names:	
OOO. Pentylone		ADB–FUBINACA)	7010
(Other names: bk-		ZZZ. methyl 2-(1-	
MBDP; 1-(1,3-		(cyclohexylmethyl)-	
benzodioxol-5-yl)-2-		1H-indole-	
(methylamino)pentan-		3-carboxamido)-3,3-	
1-one)	7542	dimethylbutanoate	
PPP. Naphyrone		(Other names:	
(Other names:		MDMB-CHMICA,	50.40
naphthylpyrovalerone; 1-		MMB-CHMINACA)	7042
(naphthalen-2-yl)-2-		AAAA. methyl 2-(1-(4-	
(pyrrolidin-1-yl)pentan-		fluorobenzyl)-	
1-one)	1258	1H-indazole-	
QQQ. alpha-pyrrolidinobutio-		3-carboxamido)-3,3-	
phenone		dimethylbutanoate (Other names:	
(Other names: α-PBP;		MDMB–FUBINACA)	7020
1-phenyl-2-(pyrrolidin-	== 4.0	BBBB. methyl 2-(1-(4-	7020
1-yl)butan-1-one)	7546	fluorobenzyl)-1H-	
RRR. N-(1-amino-3-methyl-1-		indazole-	
oxobutan-2-yl)-1-		3-carboxamido)-3-	
(cyclohexylmethyl)-		methylbutanoate	
1H-indazole-		(Other names:	
3-carboxamide		FUB–AMB, MMB–	
(Other names: AB-	5001	FUBINACA, AMB–	
CHMINACA)	7031	FUBINACA)	(7021)
SSS. N-(1-amino-3-methyl-1-		CCCC. 1-(1,3-benzodioxol-	
oxobutan-2-yl)-1-pentyl-		5-yl)-2-(ethylamino)	
1Hindazole-3-carboxamide		propan-1-one	
(Other names:	7023	(ethylone)	7547
AB-PINACA)	7023	DDDD. Naphthalen-1-yl 1-	
TTT. [1-(5-fluoropentyl)-		(5-fluoropentyl)-1H-	
1H-indazol-3-yl](naphthalen- 1-yl)methanone		indole-3-carboxylate	
(Other names:		(Other names:	
THJ-2201)	7024	NM2201; CBL2201)	7221
UUU. N-(1-amino-3,3-dimethyl-	7024	EEEE. N-(1-amino-3-methyl-	
1-oxobutan-2-yl)-1-		1-oxobutan-2-yl)-1-	
(cyclohexylmethyl)-		(5-fluoropentyl)-1H-	
1H-indazole-3-carboxamide		indazole-3-carboxamide	
(Other names: MAB-		(Other name: 5F-AB-	
CHMINACA;		PINACA)	7025
ADB-CHMINACA)	7032	FFFF. 1-(4-cyanobutyl)-N-(2-	
VVV. methyl 2-(1-(5-	,002	phenylpropan-2-yl)-1H-	
fluoropentyl)-		indazole-3-carboxamide	
1H-indazole-3-		(Other names: 4-CN- CUMYLBUTINACA;	
carboxamido)-3,3-		4-cyano-CUMYL-	
dimethylbutanoate		BUTINACA; 4-CN-	
(Other names:		CUMYLBINACA;	
5F–ADB; 5F–MDMB–		CUMYL-4CNBINACA;	
PINACA)	7034	SGT-78)	7089
WWW. methyl 2-(1-(5-		GGGG. methyl 2-(1-	,000
fluoropentyl)-		(cyclohexylmethyl)-1H-	
1H-indazole-3-		indole-3-carboxamido)-3-	
carboxamido)-3-		methylbutanoate	
methylbutanoate		(Other names: MMB-	
(Other names: 5F–		CHMICA; AMB-	
AMB)	7033	CHMICA)	7044
XXX. N-(adamantan-1-yl)-1-(5-		HHHH. 1-(5-fluoropentyl)-N-	
fluoropentyl)-1H-indazole-		(2-phenylpropan-2-yl)-	
3-carboxamide		1H-pyrrolo[2,3-b]	
(Other names: 5F–		pyridine-3-carboxamide	
APINACA, 5F–		(Other name: 5F-	
AKB48)	7049	CUMYL-P7AICA)	7085

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*			
IIII. N-ethylpentylone (Other		(I) Any compound structurally de	
names: ephylone, 1-(1,3-		3-(1-naphthoyl)indole or 1Hindol-3-yl-(1-napht	
benzodioxol-5-yl)-2-		by substitution at the nitrogen atom of the in	
(ethylamino)-pentan-1-		alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cy	
one)	7543	1-(N-methyl-2-piperidinyl)methyl or 2-(4-morp	
JJJJ. methyl 2-(1-(4-fluorobutyl)-		group, whether or not further substituted in the i	
1H-indazole-3-		any extent, whether or not substituted in the nap	ohthyl ring to
carboxamido)-3,		any extent. Including, but not limited to:	
3-dimethylbutanoate		(a) AM2201, or 1-(5-	
(4F–MDMB–BINACA,		fluoropentyl)-3-	
4F-MDMB-		(1-naphthoyl)indole	7201
BUTINACA)	7043	(b) JWH-007, or 1-	
KKKK. 1-(4-methoxyphenyl)-N-		pentyl-2-methyl-	
methylpropan-2-amine		3-(1-naphthoyl)indole	
(Other names: para-		(c) JWH-015, or 1-propyl-	
methoxymethamphetamine,	10.45	2-methyl-3-(1-	
PMMA)	1245	naphthoyl)indole	
LLLL. ethyl 2-(1-(5-fluoropentyl)-		(d) JWH-018, or 1-pentyl-	
1H-indazole-3-carboxamido)-3,3-		3-(1-naphthoyl)indole	7118
dimethylbutanoate	-	(e) JWH-019, or 1-hexyl-	
(other name: 5F-EDMB-PINACA)	7036	3-(1-naphthoyl)indole	7019
MMMM. methyl 2-(1-(5-fluoropentyl)-		(f) JWH-073, or 1-butyl-	
1 <i>H</i> -indole-3-carboxamido)-3,3-		3-(1-naphthoyl)indole	7173
Dimethylbutanoate (other names:	7041	(g) JWH-081, or 1-pentyl-	
5F-MDMB-PICA; 5F-MDMB-2201)	7041	3-(4-methoxy-1-	
NNNN. N-(adamantan-1-yl)-1-(4-		naphthoyl)indole	7081
fluorobenzyl)-1 <i>H</i> -indazole-3-		(h) JWH-098, or 1-pentyl-	
carboxamide (other names:		2-methyl-3-(4-	
FUB-AKB48; FUB-APINACA;	7047	methoxy-1-	
AKB48 N-(4-FLUOROBENZYL))	7047	naphthoyl)indole	
0000. 1-(5-fluoropentyl)-N-(2-		(i) JWH-122, or 1-pentyl-	
phenylpropan-2-yl)-1 <i>H</i> -		3-(4-methyl-1-	
indazole-3-carboxamide (other names:	7083	naphthoyl)indole	7122
5F-CUMYL-PINACA; SGT-25) PPPP. (1-(4-fluorobenzyl)-1 <i>H-</i>	7085	(j) JWH-164, or 1-pentyl-	
indol-3-yl)(2,2,3,3-tetramethylcyclopropy	4)	3-(7-methoxy-1-	
methanone (other name: FUB-144)	7014	naphthoyl)indole	
QQQQ. <i>N</i> -Ethylhexedrone (Other names:	7014	(k) JWH-200, or 1-(2-(4-	
α-ethylaminohexanophenone; 2-(ethylam	nino)	(morpholinyl)ethyl))-3-	
-1-phenylhexan-1-one)	7246	(1-naphthoyl)indole	7200
RRRR. <i>alpha</i> -Pyrrolidinohexanophenone	7240	(l) JWH-210, or 1-pentyl-	/200
(Othernames:α-PHP;α-pyrrolidinohexano	nhenone	3-(4-ethyl-1-	
1-phenyl-2-(pyrrolidin-1-yl)hexan-1-one)	7544	naphthoyl)indole	
SSSS. 4-Methyl- <i>alpha</i> -ethylaminopentioph		(m) JWH-398, or 1-pentyl-	
(Other names: 4-MEAP; 2-(ethylamino)-1	chone	3-(4-chloro-1-	
-(4-methylphenyl)pentan-1-one)	7245	naphthoyl)indole	7398
TTTT. 4'-Methyl-alpha-pyrrolidinohexioph		(II) Any compound structurally de	
(Other names: MPHP; 4'-methyl-		3-(1-naphthoyl)pyrrole by substitution at the nitro	
alpha-pyrrolidinohexanophenone; 1-(4-meth	vlphenvl)-	the pyrrole ring by alkyl, haloalkyl, alkenyl, cyclo	
2-(pyrrolidin-1-yl)hexan-1-one)	7446	cycloalkylethyl, 1-(N-methyl-2- piperidinyl)	
UUUU. alpha-Pyrrolidinoheptaphenone	,	2-(4-morpholinyl)ethyl group, whether or	
(Other names: PV8; 1-phenyl-2-		substituted in the pyrrole ring to any extent, where the pyrrole ring to any extent, where the pyrrole ring to any extent where the pyrrole ring to any extent the pyrrole ring to any ext	hether or not
(pyrrolidin-1-yl)heptan-1-one)	7548	substituted in the naphthyl ring to any extent;	
VVVV. 4'-Chloro-alpha-pyrrolidinovaleropl		(III) Any compound structurally de	erived from
(Other names: 4-chloro-α-PVP; 4'-chloro-		1-(1-naphthylmethyl)indene by substitution at the	
pyrrolidinopentiophenone; 1-(4-chloroph		the indene ring by alkyl, haloalkyl, alkenyl, cyclo	
(pyrrolidin-	· / -	cycloalkylethyl, 1-(N-methyl-2- piperidinyl)	
1-yl) pentan-1-one)	7443	2-(4-morpholinyl)ethyl group, whether or	
WWWW. 2-(ethylamino)-2-	_	substituted in the indene ring to any extent, wh	
(3-methoxyphenyl)cyclohexan-1-one		substituted in the naphthyl ring to any extent;	
(methoxetamine, MXE)	7286	(IV) Any compound structurally de	erived from
[LLLL.]XXXX. Synthetic cannabinoids:	Unless	3-phenylacetylindole by substitution at th	
ecifically exempted or unless listed in another sch		atom of the indole ring with alkyl, haloal	
stants for a state of a state of a second transition which	- contain c	avalaalludmathul avalaalludathul 1 (Nmathul )	

specifically exempted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, 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: (IV) Any compound structurally derived from 3-phenylacetylindole by substitution at the nitrogen atom of the indole ring with alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2- piperidinyl) methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent, whether or not substituted in the phenyl ring to any extent. Including, but not limited to:

<ul> <li>(a) JWH-201, or 1-pentyl- 3-(4- methoxyphenylacetyl)indole</li> <li>(b) JWH-203, or 1-pentyl- 3-(2-</li> </ul>	
chloropheny-	
lacetyl)indole	7203
(c) JWH-250, or 1-pentyl-	
3-(2-methoxypheny-	
lacetyl)indole	6250
(d) JWH-251, or 1-pentyl-	
3-(2-	
methylphenylacetyl)indole	
(e) RCS-8, or 1-(2-	
cyclohexylethyl)-3-(2-	
methoxypheny-	
lacetyl)indole	7008

(V) Any compound structurally derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the 5-position of the phenolic ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl) methyl or 2-(4-morpholinyl)ethyl group, whether or not substituted in the cyclohexyl ring to any extent. Including, but not limited to:

(a) CP 47,497 & homologues, or 2-[(1R,3S)-3hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol), where side chain n=5, and homologues where

side chain n-4, 6, or 7 7297, 7298 (VI) Any compound containing a 3- (benzoyl)indole structure with substitution at the nitrogen atom of the indole

ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl or 2-(4-morpholinyl)ethyl group, whether or not further substituted in the indole ring to any extent and whether or not substituted in the phenyl ring to any extent. Including, but not limited to:

(a) AM-694, or 1-(5-	
fluoropentyl)-3-(2-	
iodobenzoyl)indole	7694
(b) RCS-4, or 1-pentyl-3-(4-	
methoxybenzoyl)indole	
(SR-19 and RCS-4)	7104

(VII) CP 50,556-1, or [(6S,6aR,9R,10aR)-9-hydroxy-6methyl-3-[(2R)-5-phenylpentan-2-yl]oxy-5,6,6a,7,8,9,10,10aoctahydrophenanthridin- 1-yl] acetate;

(VIII) HU-210, or (6aR,10aR)-9- (hydroxymethyl)-6,6dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c] chromen-1-ol;

(IX) HU-211, Dexanabinol,(6aS,10aS)-9or (hydroxymethyl)- 6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol;

(X) Dimethylheptylpyran, or DMHP.

5. Depressants. Unless specifically excepted 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, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

A. Gamma-hydroxybutyric acid and other names gamma-hydroxybutyrate; 4-hydroxybutyrate: GHB; 4-hydroxybutonic acid; sodium oxybate; sodium oxybutryrate 2010

B. Mecloqualone	2572
C. Methaqualone	2565

C. Methaqualone

6. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, 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: A. Aminorex 1585

Some trade or other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine;

B. N-benzylpiperazine (some	
other names: BZP, 1-	
benzylpiperzaine)	7493
C. Cathinone (Some trade or	
other names: 2-amino-1-	
phenyl-1-propanone,	
alphaaminopropiophenone,	
2-aminopropiophenone and	
norephedrone)	1235
D. 4,4'-Dimethylaminorex	
(4,4'-DMAR; 4,5-dihydro-4-	
methyl-5-(4-methylphenyl)-2-	
oxazolamine; 4-methyl-5-(4-	
methylphenyl)-4,5-dihydro-	
1,3-oxazol-2-amine)	1595
E. Fenethylline	1503
F. Methcathinone	1237

Some trade or other names: 2-(methylamino)-propiophenone; alpha-(methylamino) propiophenone; 2-(methylamino)-1alpha-N-methylaminopropiophenone; phenylpropan-1-one; ephedrone; monomethylpropion; N-methylcathinone; methylcathinine; AL-464; AL-422; AL-463 and URI 432;

G. 4-methoxymethcathinone	
H. cis-4-methylaminorex	
(cis-4,5-dihydro-4-methyl-	
5-phenyl-2-oxazolamine)	1590
I. 4-Methyl-alpha-	
pyrrolidinobutiophenone,	
or MPBP	
J. N-ethylamphetamine	1475

K. N,N-dimethylamphetamine 1480 (some other names: N,N-alpha-trimethylbenzeneethanamine;

N,N-alpha-trimethylphenethylamine) 7. A temporary listing of substances subject to emergency

scheduling under federal law shall include any material, compound, mixture, or preparation which contains any quantity of the following substances:

A. Fentanyl-related substances,

their isomers, esters, ethers,

salts, and salts of isomers,	
esters, and ethers.	9850
	-

(I) 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 (1) or more of the following modifications:

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

(b) Substitution in or on the phenethyl group with alkyl, alkenyl, alkoxyl, hydroxyl, halo, haloalkyl, amino, or nitro groups;

(c) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl, ester, ether, hydroxyl, halo, haloalkyl, amino, or nitro groups;

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

(e) Replacement of the N-propionyl group by another acyl group.

47, INO. 20		
[B. ethyl 2-(1-(5-fluoropentyl)-		names: MPHP; 4'-methyl-
1H-indazole-3-carboxamido)-		alpha-pyrrolidinohexano-
,		
3,3-dimethylbutanoate, its		phenone; 1-(4-methylphenyl)-
optical, positional, and		2-(pyrrolidin-1-yl)hexan-1-
geometric isomers, salts, and		one) 7446
salts of isomers (trivial		K. alpha-Pyrrolidinohepta-
name: 5F-EDMB-		phenone, its optical,
PINACA)	7036	positional, and geometric
C. methyl 2-(1-(5-fluoropentyl)-		isomers, salts, and salts of
1H-indole-3-carboxamido)-3,3-		isomers (Other names:
dimethylbutanoate, its optical,		•
positional, and geometric		PV8; 1-phenyl-2-
		(pyrrolidin-1-yl)heptan-1-
isomers, salts, and salts of		one) 7548
isomers (trivial name:		L. 4'-Chloro-alpha-
5F-MDMB-PICA)	7041	pyrrolidinovalerophenone,
D. N-(adamantan-1-yl)-1-(4-		its optical, positional, and
fluorobenzyl)-1H-indazole-3-		geometric isomers, salts,
carboxamide, its optical,		
positional, and geometric		and salts of isomers (Other
isomers, salts, and salts		names: 4-chloro-α-PVP; 4'-
of isomers (trivial names:		chloro-alpha-
FUB-AKB48; FUB-		pyrrolidinopentiophenone; 1-
		(4-chlorophenyl)-2-(pyrrolidin-
APINACA; AKB48 N-(4-		1-vl) pentan-1-one) 7443
FLUOROBENZYL))	7047	M. N,N-diethyl-2-(2-(4
E. 1-(5-fluoropentyl)-N-(2-		
phenylpropan-2-yl)-1H-		isopropoxybenzyl)-5-nitro-
indazole-3-carboxamide, its		1H-benzimidazol-1-yl)ethan-
optical, positional, and		1-amine, its isomers,
geometric isomers, salts, and		esters, ethers, salts, and salts
salts of isomers (trivial names:		of isomers, esters, and ethers
5F-CUMYL-PINACA;		(Other names: isotonitazene;
	7083	N,N-diethyl-2-[[4- (1-
SGT-25)	7083	
F. (1-(4-fluorobenzyl)-1H-		methylethoxy)
indol-3-yl)(2,2,3,3-		phenyl]methyl]-5-nitro-1H-
tetramethylcyclopropyl)		benzimidazole-1-
methanone, its optical,		ethanamine) 9614]
positional, and geometric		[N.]B. 1-(1-(4-bromophenyl)
isomers, salts, and salts of		ethyl)piperidin-4-yl)-1, 3-
isomers (trivial name:		dihydro-2H-benzo[d]imidazol-
FUB-144)	7014	
G. N-Ethylhexedrone, its optical,	1011	2-one, its isomers, esters,
		ethers, salts, and salts of
positional, and geometric		isomers, esters, and ethers
isomers, salts, and salts of		(Other names: brorphine;
isomers (Other name: 2-		1-[1-[1-(4-bromophenyl)ethyl]-4-
(ethylamino)-1-		piperidinyl]-1,3-dihydro-
phenylhexan-1-one)	7246	2H-benzimidazol-2-one) 9098
H. alpha-Pyrrolidinohexano-		
phenone, its optical, positional,		C. 2-(2-(4-butoxybenzyl)-5-nitro-1H-
and geometric isomers, salts,		benzimidazol-1-yl)-N, N-diethylethan-1-amine,
and salts of isomers (Other		its isomers, esters, ethers, salts,
		and salts of isomers,
names: α-PHP; alpha-		esters and ethers
pyrrolidinohexiophenone; 1-		(Other name: Butonitazene) 9751
phenyl-2-(pyrrolidin-1-		D. 2-(2-(4-ethoxybenzyl)-1 <i>H</i> -benzimidazol-1-yl)- <i>N</i> ,
yl)hexan-1-one)	7544	
I. 4-Methyl-alpha-		N-diethylethan-1-amine, its isomers, esters, ethers,
ethylaminopentiophenone,		salts, and salts of isomers,
its optical, positional, and		esters and ethers
geometric isomers, salts, and		(Other names: Etodesnitazene; etazene) 9765
salts of isomers (Other		E. N, N-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-
		1H-benzimidazol-1-yl)ethan-1-amine, its isomers,
names: 4-MEAP; 2-		esters, ethers, salts, and
(ethylamino)-1-		
(4-methylphenyl)pentan-1-		salts of isomers, esters
one)	7245	and ethers (Other name: Flunitazene) 9756
J. 4'-Methyl-alpha-		F. N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-
pyrrolidinohexiophenone,		benzimidazol-1-yl)ethan-1-amine, its
its optical, positional, and		isomers, esters, ethers, salts,
		and salts of isomers, esters and ethers
geometric isomers, salts, and		
salts of isomers (Other		(Other name: Metodesnitazene) 9764

## **EMERGENCY RULES**

G. <i>N</i> , <i>N</i> -diethyl-2-(2-(4-methoxybenzyl)-5-nitro benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers	)-1 <i>H</i> -
(Other name: Metonitazene)	9757
H. 2-(4-ethoxybenzyl)-5-nitro-1-	
(2-(pyrrolidin-1-yl)ethyl) -1H-benzimidazole	,
its isomers, esters, ethers, salts,	
and salts of isomers, esters and ethers	
(Other names: N-pyrrolidino	
etonitazene; etonitazepyne)	9758
I. N, N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)	
-1H-benzimidazol-1-yl)ethan-1-amine,	
its isomers, esters, ethers, salts, and salts of is	omers,
esters and ethers	
(Other name: Protonitazene)	9759

8. Khat, to include all parts of the plant presently classified botanically as catha edulis, whether growing or not; the seeds thereof; any extract from any part of such plant; and every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seed, or extracts. 7032

(B) Schedule II shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the Controlled Substances Code Number set forth opposite it.

1. Substances, vegetable origin, or chemical synthesis. Unless specifically excepted or unless listed in another schedule, Schedule II shall include any of the following substances whether produced directly or indirectly by extraction from substances of vegetable origin or independently by means of chemical synthesis or by a combination of extraction and chemical synthesis:

A. Opium and opiate; and any salt, compound, derivative, or preparation of opium or opiate, excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, nalmefene, naloxegol, naloxone, and naltrexone and their respective salts, but including the following:

respective suits, suchieraaning the rono thing.	
(I) Raw opium	9600
(II) Opium extracts	9610
(III) Opium fluid	9620
(IV) Powdered opium	9639
(V) Granulated opium	9640
(VI) Tincture of opium	9630
(VII) Codeine	9050
(VIII) Dihydroetorphine	9334
(IX) Ethylmorphine	9190
(X) Etorphine hydrochloride	9059
(XI) Hydrocodone	9193
(XII) Hydromorphone	9150
(XIII) Metopon	9260
(XIV) Morphine	9300
(XV) Oripavine	9330
(XVI) Oxycodone	9143
(XVII) Oxymorphone	9652
(XVIII) Thebaine	9333

B. Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in subparagraph (1)(B)1.A. of this rule shall be included in Schedule II, except that these substances shall not include the isoquinoline alkaloids of opium;

C. Opium poppy and poppy

straw

D. Coca leaves (9040) and any salt, compound, derivative, or preparation of coca leaves (including cocaine (9041) and ecgonine (9180) and their salts, isomers, derivatives, and salts of isomers and derivatives), and any salt, compound, derivative, or preparation thereof which is chemically

9650

equivalent or identical with any of these substances, except that the substances shall not include:

(I) Decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine; or

(II) Ioflupane;

E. Concentrate of poppy straw (the crude extract of poppy straw in either liquid, solid, or powder form which contains the phenanthrene

alkaloids of the opium poppy) 9670 2. Opiates. Unless specifically excepted or unless in another schedule any of the following opiates, including its isomers, esters, ethers, salts, and salts of isomers, esters, and ethers whenever the existence of such isomers, esters, ethers, and salts is possible within the specific chemical designation, dextrorphan, and levopropoxyphene excepted:

A. Alfentanil	9737
B. Alphaprodine	9010
C. Anileridine	9020
D. Bezitramide	9800
E. Bulk Dextropropoxyphene	
(Non-dosage Forms)	9273
F. Carfentanil	9743
G. Dihydrocodeine	9120
H. Diphenoxylate	9170
I. Fentanyl	9801
J. Isomethadone	9226
K. Levo-alphacetylmethadol	
Some other names: levo-alphaacetylmethadol, levome	ethadyl
acetate, LAAM	9648
L. Levomethorphan	9210
M. Levorphanol 9220	
N. Metazocine	9240
O. Methadone	9250
P. Methadone-Intermediate,	
4-cyano-2-dimethylamino-	
4,4-diphenyl butane	9254
Q. Moramide-Intermediate, 2-	
methyl-3-morpholino-1,	
1-diphenylpropane-carboxylic	
acid	9802
R. Oliceridine (N-[(3-	
methoxythiophen-2-yl)	
methyl] ({2-[(9R)-9-(pyridin-	
2-yl)-6-oxaspiro	
[4.5]decan-9-	
yl]ethyl})amine fumarate)	9245
S. Pethidine (Meperidine)	9230
T. Pethidine-Intermediate-A,	
4-cyano-1-methyl-4-	
phenylpiperidine	9232
U. Pethidine-Intermediate-B,	
ethyl-4-phenylpiperidine-4-	
carboxylate	9233
V. Pethidine-Intermediate-C, 1-	
methyl-4-phenylpiperidine-	
4-carboxylic acid	9234
W. Phenazocine	9715
X. Piminodine	9730
Y. Racemethorphan	9732
Z. Racemorphan	9733
AA. Remifentanil	9739
BB. Sufentanil	9740
CC. Tapentadol	9780
DD. Thiafentanil	9729

3. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following

substances having a stimulant effect on the central nervous system:

A. Amphetamine, its salts,	
optical isomers, and salts	
of its optical isomers	1100
B. Lisdexamfetamine, its salts,	
isomers, and salts of its	
isomers	1205
C. Methamphetamine, its salts,	
isomers, and salts of its	
isomers	1105
D. Phenmetrazine and its salts	1631
E. Methylphenidate	1724

4. Depressants. Unless specifically excepted 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, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

A. Amobarbital	2125
B. Glutethimide	2550
C. Pentobarbital	2270
D. Phencyclidine	7471
E. Secobarbital	2315
5. Hallucinogenic substances:	
A. Nabilone	7379

Another name for nabilone: (±)trans-3-(1, 1dimethylheptyl)-6, 6a,7,8,10,10a-hexahydro- 1-hydroxy-6, 6-dimethyl-9H-dibenzo(b,d) pyran-9-one.

B. Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] in an oral solution in a drug product approved for marketing by the United States Food and Drug Administration. (7365)

6. Immediate precursors. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances:

A. Immediate precursor to amphetamine and methamphetamine:

(I) Phenylacetone 8501 Some trade or other names: phenyl-2-propanone; P2P; benzyl methyl ketone; methyl benzyl ketone;

,	
B. Immediate precursors to phencyclidine (PCP):	
(I) 1-phenylcyclohexylamine	7460
(II) 1-piperidinocyclo-	
hexanecarbonitrile (PCC)	8603
C. Immediate precursor to fentanyl:	
(I) 4-anilino-N-phenethyl-4-	
piperidine (ANPP)	8333
(II) N-phenyl-N-(piperidin-	
4-yl)propionamide (norfentanyl)	8366

7. Any material, compound, mixture, or preparation which contains any quantity of the following alkyl nitrites:

A. Amyl nitrite;

B. Butyl nitrite.

(C) Schedule III shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this section. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.

1. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, 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 (whether optical, position, or geometric), and salts of such isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation: A. Those compounds, mixtures, or preparations in dosage unit form containing any stimulant substances listed in Schedule II which compounds, mixtures, or preparations were listed on August 25, 1971, as excepted compounds under 21 CFR 308.32 and any other drug of the quantitive composition shown in that list for those drugs or which is the same except that it contains a lesser quantity of controlled substances 1405

trolled substances	1405
B. Benzphetamine	1228
C. Chlorphentermine	1645
D. Clortermine	1647
E. Phendimetrazine	1615

2. Depressants. Unless specifically excepted 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:

	A. Any	compound,	mixture,	or	preparation
contai	ning –				
	(I) Amo	obarbital			2126
(II) Secobarbital					2316

(III) Pentobarbital 2271 or any salt thereof and one (1) or more other active medicinal ingredients which are not listed in any schedule:

lients which are not listed in any schedule;	
B. Any suppository dosage form containing –	
(I) Amobarbital	2126
(II) Secobarbital	2316
(III) Pentobarbital	2271

(III) Pentobarbital 2271 or any salt of any of these drugs and approved by the Food and Drug Administration for marketing only as a suppository;

C. Any substance which contains any
quantity of a derivative of barbituric acid
or any salt thereof

ny salt thereof	2100
D. Chlorhexadol	2510
E. Embutramide	2020

F. Any drug product containing gamma hydroxybutyric acid, including its salts, isomers, and salts of isomer, for which an application is approved under section 505 of the Federal Food, Drug, and Cosmetic Act; 2012

-	5.	
	G. Ketamine, its salts, isomer,	
	and salts of isomers (some	
	other names for ketamine:	
	(±)-2-(2-chlorophenyl)-2-	
	(methylamino)-	
	cyclohexanone)	7285
	H. Lysergic acid	7300
	I. Lysergic acid amide	7310
	J. Methyprylon	2575
	K. Perampanel, and its salts,	
	isomers, and salts of	
	isomers	2261
	L. Sulfondiethylmethane	2600
	M. Sulfonethylmethane	2605
	N. Sulfonmethane	2610
	O. Tiletamine and zolazepam	
	or any salt thereof	7295
	-	

Some trade or other names for a tiletaminezolazepam 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- trimethylpyrazolo-(3,4-e) (1,4)-diazepin-7(1H)-one, flupyrazapon.

3. Nalorphine

9400

4. Narcotics drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof: one)

#### A. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of opium 9803

B. Not more than one and eight tenths grams (1.8gm) of codeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9804

C. Not more than one and eight tenths grams (1.8gm) of dihydrocodeine per one hundred milliliters (100 mL) or not more than ninety milligrams (90 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9807

D. Not more than three hundred milligrams (300 mg) of ethylmorphine per one hundred milliliters (100 mL) or not more than fifteen milligrams (15 mg) per dosage unit, with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9808

E. Not more than five hundred milligrams (500 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm) or not more than twenty-five milligrams (25 mg) per dosage unit, with one (1) or more active nonnarcotic ingredients in recognized therapeutic amounts 9809

F. Not more than fifty milligrams (50 mg) of morphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm), with one (1) or more active, nonnarcotic ingredients in recognized therapeutic amounts 9810

5. Any material, compound, mixture, or preparation containing any of the following narcotic drugs or their salts, as set forth below:

A. Buprenorphine

9064

6. Anabolic steroids. Unless specially excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts of isomers is possible within the specific chemical designation. DEA has assigned code 4000 for all anabolic steroids. Anabolic steroids. Any drug or hormonal substance, chemically and pharmacologically related to testosterone (other than estrogens, progestins, corticosteroids, and dehydroepiandrosterone) that promotes muscle growth, except an anabolic steroid which is expressly intended for administration through implants to cattle or other nonhuman species and which has been approved by the Secretary of Health and Human Services for that administration. If any person prescribes, dispenses, or distributes such steroid for human use, such person shall be considered to have prescribed, dispensed, or distributed an anabolic steroid within the meaning of this paragraph. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any quantity of the following substances, including its salts, esters, and ethers:

- A.  $3\beta$ ,  $17\beta$ -dihydroxy- $5\alpha$ -androstane
- B. 3α,17β-dihydroxy-5α-androstane
- C. 5q-androstan-3,17-dione
- D. 1-androstenediol
  - (3β,17β-dihydroxy-5α-androst-1-ene)
- E. 1-androstenediol  $(3\alpha, 17\beta$ -dihydroxy- $5\alpha$ -androst-1-ene)
- F. 4-androstenediol (3β,17β-dihydroxy-androst-4-ene)
- G. 5-androstenediol ( $3\beta$ ,  $17\beta$ -dihydroxy-androst-5-ene)
- H. 1-androstenedione ([5α]-androst-1-en-3,17-dione)
- I. 4-androstenedione (androst-4-en-3,17-dione)
- J. 5-androstenedione (androst-5-en-3,17-dione)
- K. Bolasterone (7α,17α-dimethyl-17β-
- hydroxyandrost-4-en-3- one)
- L. Boldenone (17β-hydroxyandrost-1, 4-diene-3-one)

- M. Boldione (androstra-1,4-diene-3,17-dione)
- N. Calusterone (7β,17α-dimethyl-17β-
- hydroxyandrost-4-en-3-one)
- O. Clostebol (4-chloro-17β-hydroxyandrost-4-en-3-one)
- P. Dehydrochloromethyltestosterone  $(4-chloro-17\beta-hydroxy-17\alpha-methyl-androst-1,$ 4-dien-3-one)
- Q. Desoxymethyltestosterone  $(17\alpha$ -methyl- $5\alpha$ -androst-2-en- $17\beta$ -ol) (a.k.a. madol)
- R.  $\Delta$ 1-dihydrotestosterone (a.k.a.'1-testosterone')(17β- hydroxy-5α-androst-1-en-3-
- S. 4-dihydrotestosterone (17β-hydroxy-androstan-3-one)
- T. Drostanolone (17β-hydroxy-2α-methyl-5α-androstan-3-one)
- U. Ethylestrenol (17α-ethyl-17β-hydroxyestr-4-ene)
- V. Fluoxymesterone (9-fluoro-17α-methyl-11β, 17β-dihydroxyandrost-4-en-3-one)
- W. Formebulone (Formebolone) (2-formyl-17αmethyl-11 $\alpha$ ,17 $\beta$ -dihydroxyandrost-1,4-dien-3-one)
- X. Furazabol (17α-methyl-17β-hydroxyandrostano [2,3-c]-furazan)
- Y. 13β-ethyl-17β-hydroxygon-4-en-3-one
- Z. 4-hydroxytestosterone (4,17β-dihydroxy-androst-4-en-3-one)
- AA. 4-hydroxy-19-nortestosterone (4,17β-dihydroxy-estr-4-en-3-one)
- BB. Mestanolone
- $(17\alpha$ -methyl-17 $\beta$ -hydroxy-5 $\alpha$ -androstan-3-one) CC. Mesterolone
- $(1\alpha$ -methyl-17 $\beta$ -hydroxy- $(5\alpha)$ -androstan-3-one) DD. Methandienone

(17α-methyl-17β-hydroxyandrost-1,4-dien-3-one)

- EE. Methandriol (17α-methyl-3β,17β-
- dihydroxyandrost-5-ene)

FF. Methasterone (2α,17α-dimethyl-5α-androstan-17β-ol-3-one)

GG. Methenolone (1-methyl-17β-hydroxy-5α-androst-1en-3-one)

HH.  $17\alpha$ -methyl- $3\beta$ , $17\beta$ -dihydroxy- $5\alpha$ -androstane

II.  $17\alpha$ -methyl- $3\alpha$ ,  $17\beta$ -dihydroxy- $5\alpha$ -androstane

JJ. 17α-methyl-3β,17β-dihydroxyandrost-4-ene

KK. 17α-methyl-4-hydroxynandrolone(17α-methyl-4hydroxy-17β-hydroxyestr-4- en-3-one)

LL. Methyldienolone (17α-methyl-17β-hydroxyestra-4,9(10)-dien-3-one)

MM. Methyltrienolone (17α-methyl-17β-hydroxyestra-4,9,11-trien-3-one)

NN. Methyltestosterone (17α-methyl-17hydroxyandrost-4-en-3-one)

OO. Mibolerone (7α,17α-dimethyl-17β-hydroxyestr-4-en-3-one)

PP. 178-methyl- $\Delta$ 1-dihydrotestosterone (17β-hydroxy- $17\alpha$ -methyl- $5\alpha$ -androst-1-en-3-one) (a.k.a. 17-α-methyl-1testosterone)

- QQ. Nandrolone (17 $\beta$ -hydroxyestr-4-ene-3-one)
- RR. 19-nor-4-androstenediol (3β,17β-dihydroxyestr-4-ene)
- SS. 19-nor-4-andro stenediol (3α,17β-dihydroxyestr-4-ene)
- TT. 19-nor-4,9(10)-androstadienedione (estra-4,9(10)-
- diene-3,17-dione)
  - UU. 19-nor-5-androstenediol (3β,17β-dihydroxyestr-5-ene)
  - VV. 19-nor-5-androstenediol (3α,17β-dihydroxyestr-5-ene)
  - WW. 19-nor-4-androstenedione (estr-4-en-3,17-dione) XX. 19-nor-5-androstenedione (estr-5-en-3,17-dione)
- YY. Norbolethone (13β,17α-diethyl-17β-hydroxygon-4en-3-one)

ZZ. Norclostebol (4-chloro-17β-hydroxyestr-4-en-3-one)

AAA. Norethandrolone (17 $\alpha$ -ethyl-17 $\beta$ -hydroxyestr-4-en-3-one)

BBB. Normethandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxyestr-4-en-3-one)

CCC. Oxandrolone (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-2-oxa-[5 $\alpha$ ]- androstan-3-one)

DDD. Oxymesterone (17α-methyl-4,17βdihydroxyandrost-4-en-3-one)

EEE. Oxymetholone (17α-methyl-2-hydroxymethylene-17β-hydroxy-[5α]-androstan-3-one)

FFF. Prostanozol (17 $\beta$ -hydroxy-5 $\alpha$ -androstano[3,2-c] pyrazole)

GGG. Stanolone ( $\Delta$ 1-dihydrotestosterone (a.k.a. 1-testosterone)(17 $\beta$ -hydroxy-5 $\alpha$ -androst-1-en-3-one))

HHH. Stanozolol (17 $\alpha$ -methyl-17 $\beta$ -hydroxy-[5 $\alpha$ ]-androst-2-eno[3,2-c]-pyrazole)

III. Stenbolone (17 $\beta$ -hydroxy-2-methyl-[5 $\alpha$ ]-androst-1-en-3-one)

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

KKK. Testosterone( $17\beta$ -hydroxyandrost-4-en-3-one);

LLL. Tetrahydrogestrinone (13 $\beta$ ,17 $\alpha$ -diethyl-17 $\beta$ -hydroxygon-4,9, 11-trien-3-one)

MMM. Trenbolone ( $17\beta$ -hydroxyestr-4,9,11-trien-3-one)

NNN. Any salt, ester, or isomer of a drug or substance described or listed in this subparagraph, if that salt, ester, or isomer promotes muscle growth except an anabolic steroid which is expressly intended for administration through implants to cattle or other nonhuman species and which has been approved by the Secretary of Health and Human Services for that administration.

7. Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a United States Food and Drug Administration approved drug product 7369 (Some other names for dronabinol: (6aRtrans)- 6a,7,8,10atetrahydro-6.6.9-trimethyl-3-pentyl-6H-dibenzo (b,d) pyran-1ol, or (-) -delta-9-(trans)-tetrahydrocannabinol.)

(D) Schedule IV shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical name, or brand name designated, listed in this subsection. Each drug or substance has been assigned the DEA Controlled Substances Code Number set forth opposite it.

1. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing limited quantities of any of the following narcotic drugs or any salts thereof:

A. Not more than one milligram (1 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit 9167

B. Dextropropoxyphene (alpha-(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-propionoxybutane) 9278

C. 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl) cyclohexanol, its salts, optical and geometric isomers, and salts of these isomers (including tramadol) 9752

D. Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following limited quantities of narcotic drugs or salts thereof, which shall include one (1) or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:

(I) Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

(II) Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm); or

(III) Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm).

2. Depressants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:

ers is possible within the specific chemical des	ignation:
A. Alfaxalone	2731
B. Alprazolam	2882
C. Barbital	2145
D. Brexanolone	2400
E. Bromazepam	2748
F. Camazepam	2749
G. Carisoprodol	8192
H. Chloral betaine	2460
I. Chloral hydrate	2465
J. Chlordiazepoxide	2744
K. Clobazam	2751
L. Clonazepam	2737
M. Clorazepate	2768
N. Clotiazepam	2752
O. Cloxazolam	2753
P. Daridorexant	<b>2410</b>
[P.]Q. Delorazepam	2754
[Q.]R. Diazepam	2765
[R.]S. Dichloralphenazone	2467
[S.JT. Estazolam	2407
[ <b>T</b> .] <b>U.</b> Ethchlorvynol	2540
[U.]V. Ethinamate	2545
[V.]W. Ethyl loflazepate	2758
[W.]X. Fludiazepam	2759
[X.]Y. Flunitrazepam	2763
[Y.]Z. Flurazepam	2767
[Z.]AA. Fospropofol	2138
[AA.]BB. Halazepam	2762
[BB.]CC. Haloxazolam	2771
[CC.]DD. Ketazolam	2772
[DD.]EE. Lemborexant	2245
[EE.]FF. Loprazolam	2773
[FF.]GG. Lorazepam	2885
[GG.]HH. Lormetazepam	2774
[HH.]II. Mebutamate	2800
[//.]JJ. Medazepam	2836
[JJ.]KK. Meprobamate	2820
[KK.]LL. Methohexital	2264
[LL.]MM. Methylphenobarbital	
(Mephobarbital)	2250
<i>[MM.]</i> NN. Midazolam	2884
[NN.]OO. Nimetazepam	2837
[OO.]PP. Nitrazepam	2834
[PP.]QQ. Nordiazepam	2838
[QQ.]RR. Oxazepam	2835
[RR.]SS. Oxazolam	2839
[SS.]TT. Paraldehyde	2585
[TT.]UU. Petrichloral	2591
[UU.]VV. Phenobarbital	2285
VV. WW. Pinazepam	2883
[WW.]XX. Prazepam	2764
[XX.]YY. Quazepam	2881
YY. <b>JZZ.</b> Remimazolam	2846
[ZZ.]AAA. Suvorexant	2223
[AAA.]BBB. Temazepam	2925
[BBB.]CCC. Tetrazepam	2886
[CCC.]DDD. Triazolam	2887
[DDD.]EEE. Zaleplon	2781
[EEE.]FFF. Zolpidem	2783
·	1,00

#### [FFF.]GGG. Zopiclone

3. Fenfluramine. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers (whether optical, position, or geometric), and salts of such isomers, whenever the existence of such salts, isomers, and salts of isomers is possible:

A. Fenfluramine

1670

2784

4. Lorcaserin. Any material, compound, mixture, or preparation which contains any quantity of the following substances, including its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers is possible:

A. Lorcaserin

1625

5. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, 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:

A. Cathine ((+)-	
norpseudoephedrine)	1230
B. Diethylpropion	1610
C. Fencamfamin	1760
D. Fenproporex	1575
E. Mazindol	1605
F. Mefenorex	1580
G. Modafinil	1680
H. Pemoline (including	
organometallic complexes	
and chelates thereof)	1530
I. Phentermine	1640
J. Pipradrol	1750
K. Serdexmethylphenidate	1729
L. Sibutramine	1675
M. Solriamfetol (2-amino-3-	
phenylpropyl carbamate;	
benzenepropanol, beta-	
amino-, carbamate (ester))	1650
N. SPA (-)-1-dimethylamino-	
1,2-diphenylethane	1635

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

A. Pentazocine	9709
B. Butorphanol (including its	
optical isomers)	9720
C. Eluxadoline (5-[[[(2S)-2-	
amino-3-[4-aminocarbonyl)-	
2,6-dimethylphenyl]-1-	
oxopropyl] [(1S)-1-(4-phenyl-	
1 H-imidazol-2-	
yl)ethyl]amino]methyl]-2-	
methoxybenzoic acid)	
(including its optical isomers)	
and its salts, isomers, and	
salts of isomers	9725

7. Ephedrine. Any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system including their salts, isomers, and salts of isomers:

A. Ephedrine or its salts, optical isomers, or salts of optical isomers as the only active medicinal ingredient or contains ephedrine or its salts, optical isomers, or salts of optical isomers and therapeutically insignificant quantities of another active medicinal ingredient.

(E) Schedule V shall consist of the drugs and other substances, by whatever official name, common or usual name, chemical

name, or brand name designated, listed in this subsection.

1. Narcotic drugs containing nonnarcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as follows, which shall include one (1) or more nonnarcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by the narcotic drug alone:

A. Not more than two hundred milligrams (200 mg) of codeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

B. Not more than one hundred milligrams (100 mg) of dihydrocodeine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

C. Not more than one hundred milligrams (100 mg) of ethylmorphine per one hundred milliliters (100 mL) or per one hundred grams (100 gm);

D. Not more than two and five-tenths milligrams (2.5 mg) of diphenoxylate and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit;

E. Not more than one hundred milligrams (100 mg) of opium per one hundred milliliters (100 mL) or per one hundred grams (100 gm); and

F. Not more than five-tenths milligram (0.5 mg) of difenoxin (DEA Drug Code No. 9168) and not less than twenty-five micrograms (25 mcg) of atropine sulfate per dosage unit.

2. Stimulants. 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 stimulant effect on the central nervous system

including its salts, isomers, and salts of isomers:

A. Pyrovalerone 1485 3. Any compound, mixture, or preparation containing any detectable quantity of pseudoephedrine or its salts or optical isomers, or salts of optical isomers or any compound, mixture, or preparation containing any detectable quantity of ephedrine or its salts or optical isomers, or salts of optical isomers if the drug preparations are starch-based solid dose forms, if such preparations are sold over the counter without a prescription. The following drug preparations containing ephedrine and pseudoephedrine are not scheduled controlled substances:

A. Drug preparations in liquid form;

B. Drug preparations that require a prescription in order to be dispensed.

4. 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:

A. Ezogabine [N-[2-amino-4(4-	
fluorobenzylamino)-phenyl]-	
carbamic acid ethyl ester]	2779
B. Ganaxolone (3α-hydroxy-3β-methyl-	
5α-pregnan-20-one)	2401
[B.]C. Lacosamide [(R)-2-	
acetoamido-N-benzyl-	
3-methoxy-propionamide]	2746
[C.]D. Pregabalin [(S)-3-	
(aminomethyl)-5-	
methylhexanoic acid]	2782
[D.]E. Brivaracetam ((25)-2-[(4R)-	
2-oxo-4-propylpyrrolidin-1-	
yl]butanamide) (also referred	
to as BRV; UCB-34714;	
Briviact)	2710

[E.]F. Lasmiditan [2,4,6- trifluoro-N-(6-(1- methylpiperidine-4- carbonyl) pyridine-2- yl-benzamide]	2790
[F.JG. Cenobamate ([(1R)-1-(2- chlorophenyl)-2-(tetrazol-	
2-yl)ethyl] carbamate; 2H-	
tetrazole-2-ethanol, alpha-(2- chlorophenyl)-, carbamate	
(ester), (alphaR)-; carbamic acid (R)-(+)-1-(2-	
chlorophenyl)-2-(2H-tetrazol-	0700
2-yl)ethyl ester)	2720

AUTHORITY: section 195.015, RSMo Supp. [2021] 2022, and section 195.195, RSMo 2016. Material found in this rule previously filed as 19 CSR 30-1.010. Original rule filed April 14, 2000, effective Nov. 30, 2000. For intervening history, please consult the Code of State Regulations. Emergency amendment filed Sept. 12, 2022, effective Oct. 3, 2022, expires March 31, 2023. A proposed amendment covering this same material is published in this issue of the Missouri Register.

PUBLIC COST: This emergency amendment will not cost state agencies or political subdivisions more than five hundred dollars (\$500) in the time the emergency is effective.

PRIVATE COST: This emergency amendment will not cost private entities more than five hundred dollars (\$500) in the time the emergency is effective.

#### Title 19 – DEPARTMENT OF HEALTH AND SENIOR SERVICES Division 30 – Division of Regulation and Licensure Chapter 20 – Hospitals

#### **EMERGENCY RULE**

# 19 CSR 30-20.144 Standards and Guidelines for Essential Caregiver Program

PURPOSE: This rule establishes the standards and guidelines regarding the essential caregiver program established under section 191.2290, RSMo.

EMERGENCY STATEMENT: The "Essential Caregiver Program Act" became law on August 28, 2022. Under this new law, the department is required to develop standards and guidelines concerning the essential caregiver program. These standards and guidelines will provide the regulatory framework hospitals must follow to ensure that their patients have access to their designated essential caregivers. The standards and quidelines developed by the department must be operational during a state of emergency declared pursuant to Chapter 44, relating to infectious, contagious, communicable, or dangerous diseases. While there is no current declaration of a state of emergency, the implementing regulations for the Essential Caregiver Program Act should be in place prior to that potential event. It is imperative that this rule become effective as close to the same time that the law becomes effective in order to ensure that a hospital patient has immediate access to his or her essential caregiver, an indispensable member of the patient's care team. As a result, the department finds a compelling governmental interest, which requires this emergency action. A proposed rule, which covers the same material, is published in this issue of the Missouri Register. The scope of this emergency rule is limited to the circumstances creating the emergency and complies with the protections extended in the **Missouri** and **United States Constitutions**. The department believes this emergency rule is fair to all interested persons and parties under the circumstances. This emergency rule was filed September 15, 2022, becomes effective September 29, 2022, and expires March 27, 2023.

(1) As used in this rule, the following terms and phrases shall mean:

(A) Department shall mean the Department of Health and Senior Services;

(B) Essential caregiver shall mean a family member, friend, guardian, or other individual selected by a hospital patient who has not been adjudged incapacitated under chapter 475, or the guardian or legal representative of the patient;

(C) Hospital shall have the same meaning assigned to it 19 CSR 30-20.011(9).

(2) Every hospital within Missouri shall develop an essential caregiver program which shall allow a patient who has not been adjudged incapacitated under chapter 475, RSMo, a patient's guardian, or a patient's legally authorized representative to designate an essential caregiver for in-person contact with the patient in accordance with the provisions of section 191.2290, RSMo, and the standards and guidelines developed by the department under this rule.

(3) The essential caregiver program shall be operable during a state of emergency declared pursuant to chapter 44, RSMo, relating to infectious, contagious, communicable, or dangerous diseases.

(4) The essential caregiver program established by the hospital shall:

(A) Allow at least two individuals per patient to be designated as essential caregivers, although the hospital may limit the in-person contact to one caregiver at a time. The caregiver shall not be required to have previously served in a caregiver capacity prior to the declared state of emergency;

(B) Include a reasonable in-person contact schedule to allow the essential caregiver to provide care to the patient for at least four (4) hours each day, including evenings, weekends, and holidays, but shall allow for twenty-four-hour in-person care as necessary and appropriate for the well-being of the patient. The essential caregiver shall be permitted to leave and return during the scheduled hours or be replaced by another essential caregiver;

(C) Include procedures to enable physical contact between the patient and the essential caregiver. The hospital may not require the essential caregiver to undergo more stringent screening, testing, hygiene, personal protective equipment, and other infection control and prevention protocols than required of hospital employees; and

(D) Specify in its protocols the criteria that the hospital will use if it determines that in-person contact by a particular essential caregiver is inconsistent with the patient's therapeutic care and treatment or is a safety risk to other patients or staff at the facility. Any limitations placed upon a particular essential caregiver shall be reviewed and documented every seven (7) days to determine if the limitations remain appropriate.

(5) A hospital shall inform, in writing, patients who have not been adjudged incapacitated under chapter 475, or guardians or legal representatives of patients, of the essential caregiver program and the process for designating an essential caregiver. Consistent with 42 CFR 482.12(h), a hospital shall inform each patient, or such patient's guardian or legal representative, where appropriate, of his or her visitation rights and right to access an essential caregiver in accordance with this rule.

(6) A hospital may restrict or revoke in-person contact by an essential caregiver who fails to follow required protocols and procedures established under section (4) of this rule.

(7) A hospital may request from the department a suspension of in-person contact by essential caregivers for a period not to exceed seven (7) days. A hospital may request from the department an extension of a suspension for more than seven (7) days, but such extension period shall not be for a period longer than seven (7) days at a time. Under the provisions of this section, a hospital shall not suspend in-person caregiver contact for more than fourteen (14) consecutive days in a twelve-month period or for more than forty-five (45) total days in a twelve-month period. Requests for a suspension of in-person contact of essential caregivers or an extension of a suspension under this section shall be submitted in writing to the department. Department determinations in response to suspension requests shall be in writing and both requests and determinations shall be made a part of the department's permanent records for the hospital.

(A) Requests for a suspension of in-person contact by essential caregivers shall contain at a minimum the following:

1. The specific reason or reasons why allowing in-person contact by essential caregivers poses a serious community health risk;

2. An explanation of the extenuating factors which may be relevant to granting a suspension to the particular requesting hospital; and

3. The length of time, not to exceed seven (7) days, the suspension is being requested.

(8) The department's written determination shall identify a suspension expiration date, if approved. The hospital may reapply for an extension of the suspension up to one (1) day prior to the expiration of the department's originally approved suspension. The department may deny a hospital's request to suspend in-person contact with essential caregivers if the department determines that such in-person contact does not pose a serious community health risk.

(9) The department shall suspend in-person contact by essential caregivers under this rule if it determines that doing so is required under federal law, including a determination that federal law requires a suspension of in-person contact by members of the patient's care team.

(10) The provisions of this rule shall not apply to those patients whose particular plan of therapeutic care and treatment necessitates restricted or otherwise limited visitation for reasons unrelated to the stated reasons for the declared state emergency.

(11) The provisions of this rule shall not be construed to require an essential caregiver to provide necessary care to a patient and a hospital shall not require an essential caregiver to provide necessary care.

AUTHORITY: sections 191.2290 and 197.080, RSMo Supp. 2022. Emergency rule filed Sept. 15, 2022, effective Sept. 29, 2022, expires March 27, 2022. A proposed rule covering this same material is published in this issue of the **Missouri Register**.

PUBLIC COST: This emergency rule will cost state agencies or political subdivisions one hundred twenty-two thousand four hundred dollars (\$122,400) in the time the emergency is effective. PRIVATE COST: This emergency rule will cost private entities four hundred thirty-eight thousand six hundred dollars (\$438,600) in the time the emergency is effective.

#### FISCAL NOTE PUBLIC COST

# I.Department Title:Department of Health and Senior ServicesDivision Title:Division 30—Division of Regulation and LicensureChapter Title:Chapter 20—Hospitals

Rule Number and Title:	19 CSR 30-20.144 Standards and Guidelines for Essential Caregiver Program.	
Type of Rulemaking:	Emergency Rulemaking	

### II. SUMMARY OF FISCAL IMPACT

Estimate of the number of entities by class which would likely be affected by the adoption of the rule:	Classification by types of the business entities which would likely be affected:	Estimate in the aggregate as to the cost of compliance with the rule by the affected entities:
36 public hospitals	Public Hospitals	\$122,400.00

### III. WORKSHEET Cost for Private Hospitals to Adopt and Implement Essential Caregiver Programs

Action	Explanation	Cost	Cost for Private Hospitals
Policy and procedure development, implementation, and training	Policy and Procedure Develoment-1FTE*8hrs=\$320 Implementation- 1FTE*2hrs=80 Training-100 FTE*1hr=\$3000	\$3400	36 public hospitals * \$3400 = \$122,400

#### IV. ASSUMPTIONS

While it is generally assumed that most hospitals have already built into their operational costs the cost of updating their individual institutional policies and procedures to reflect changes made in law, this fiscal note attempts to breakdown the individual cost of complying with §191.2290, RSMo and the proposed emergency rule. In order to comply with the provisions of the proposed emergency rule, hospitals will have to update their

visitation policies to incorporate the essential caregiver guidelines and standards established by the proposed emergency rule.

This fiscal note also assumes that a state of emergency under Chapter 44, RSMo, relating to infectious diseases, has not been declared and is not in place. Even though the Department does not expect a state of emergency to be declared or in place during the time period of this emergency rule, the Department does expect public hospitals to adopt and implement policies to be in compliance with the provisions of the emergency rule. Of course, the steps taken by public hospitals to implement policies and train personnel to be consistent with the essential caregiver emergency rule will have a fiscal impact.

The department licenses approximately 36 public hospitals. The Department estimates that each public hospital will incur approximately \$3,400 in costs in developing the policies and procedures for the implementation of this emergency rule.

#### FISCAL NOTE PRIVATE COST

I.	<b>Department Title:</b>	Department of Health and Senior Services
	<b>Division</b> Title:	<b>Division 30—Division of Regulation and Licensure</b>
	Chapter Title:	Chapter 20 — Hospitals

Rule Number and Title:	19 CSR 30-20.144 Standards and Guidelines for Essential Caregiver Program.
Type of Rulemaking:	Emergency Rulemaking

### II. SUMMARY OF FISCAL IMPACT

Estimate of the number of entities by class which would likely be affected by the adoption of the rule:	Classification by types of the business entities which would likely be affected:	Estimate in the aggregate as to the cost of compliance with the rule by the affected entities:
129	Private Hospitals	\$438,600.00 for 6 month period that a state of emergency is in effect (assumes no state of emergency has been declared or is in place)

#### III. WORKSHEET

### Cost for Private Hospitals to Adopt and Implement Essential Caregiver Programs

Action	Explanation	Cost	Cost for Private Hospitals
Policy and procedure development, implementation, and training	Policy and Procedure Develoment-1FTE*8hrs=\$320 Implementation- 1FTE*2hrs=80 Training-100 FTE*1hr=\$3000	\$3400	129 private hospitals * \$3400 = \$438,600.00

#### IV. ASSUMPTIONS

While it is generally assumed that most hospitals have already built into their operational costs the cost of updating their individual institutional policies and procedures to reflect changes made in law, this fiscal note attempts to breakdown the individual cost of complying with §191.2290, RSMo and the proposed emergency rule. In order to comply with the provisions of the proposed emergency rule, hospitals will have to update their visitation policies to incorporate the essential caregiver guidelines and standards established by the proposed emergency rule.

This fiscal note also assumes that a state of emergency under Chapter 44, RSMo, relating to infectious diseases, has not been declared and is not in place. Even though the Department does not expect a state of emergency to be declared or in place during the time period of this emergency rule, the Department does expect private hospitals to adopt and implement policies to be in compliance with the provisions of the emergency rule. Of course, the steps taken by private hospitals to implement policies and train personnel to be consistent with the essential caregiver emergency rule will have a fiscal impact.

The department licenses approximately 129 private hospitals (hospitals not owned by state or local governments). The Department estimates that each private hospital will incur approximately \$3,400 in costs in developing the policies and procedures for the implementation of this emergency rule.