§ 1308.01

AUTHORITY: 21 U.S.C. 811, 812, 871(b), 956(b), unless otherwise noted.

SOURCE: 38 FR 8254, Mar. 30, 1973, unless otherwise noted. Redesignated at 38 FR 26609, Sept. 24, 1973.

GENERAL INFORMATION

§ 1308.01 Scope of this part.

Schedules of controlled substances established by section 202 of the Act (21 U.S.C. 812) and nonnarcotic substances, chemical preparations, veterinary anabolic steroid implant products, prescription products, anabolic steroid products, and cannabis plant material and products made therefrom that contain tetrahydrocannabinols excluded pursuant to section 201 of the Act (21 U.S.C. 811), as they are changed, updated, and republished from time to time, are set forth in this part.

[81 FR 97021, Dec. 30, 2016]

§1308.02 Definitions.

Any term contained in this part shall have the definition set forth in section 102 of the Act (21 U.S.C. 802) or part 1300 of this chapter.

[62 FR 13967, Mar. 24, 1997]

§ 1308.03 Administration Controlled Substances Code Number.

(a) Each controlled substance, or basic class thereof, has been assigned an "Administration Controlled Substances Code Number" for purposes of identification of the substances or class on certain Certificates of Registration issued by the Administration pursuant to §§1301.35 of this chapter and on certain order forms issued by Administration pursuant §1305.05(d) of this chapter. Applicants for procurement and/or individual manufacturing quotas must include the appropriate code number on the application as required in §§1303.12(b) and 1303.22(a) of this chapter. Applicants for import and export permits must include the appropriate code number on the application as required in §§ 1312.12(a) and 1312.22(a) of this chapter. Authorized registrants who desire to import or export a controlled substance for which an import or export permit is not required must include the appropriate Administration Controlled Substances Code Number beneath or beside the name of each controlled substance listed on the DEA Form 236 (Controlled Substance Import/Export Declaration) which is executed for such importation or exportation as required in §§ 1312.18(c) and 1312.27(b) of this chapter.

(b) Except as stated in paragraph (a) of this section, no applicant or registrant is required to use the Administration Controlled Substances Code Number for any purpose.

[38 FR 8254, Mar. 30, 1973. Redesignated at 38 FR 26609, Sept. 24, 1973 and amended at 51 FR 15318, Apr. 23, 1986; 62 FR 13968, Mar. 24, 1997]

SCHEDULES

§ 1308.11 Schedule I.

- (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 DEA Controlled Substances Code Number set forth opposite it.
- (b) 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 (for purposes of 3-methylthiofentanyl only, the term isomer includes the optical and geometric isomers):

(1)	Acetyl-alpha-methylfentanyl	$(N-\Gamma 1-(1-\text{methvl}-2-\text{phenethvl})-4-$	
	peridinyl]-N-phenylacetamide)		9815
	cetylmethadol		9601
(3) A	cetyl fentanyl (N-(1-phenethylpipe:	ridin-4-yl)-N-phenylacetamide)	9821
(4)	Acryl fentanyl (N-(1-phenethylpi)	peridin-4-yl)-N-phenylacrylamide;	
als	o known as acryloylfentanyl)		9811
(5)	AH-7921	(3,4-dichloro-N-[(1-	
dir	nethylamino)cyclohexylmethyl]bei	nzamide)	9551

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(6) Allylprodine	
(7) Alphacetylmethadol (except levo-alphacetylmethadol also known as levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM)	3
(8) Alphameprodine	
(9) Alphamethadol	•
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	-
propaniido)piperidinė) (11) $alpha$ -Methylthiofentanyl (N -[1-methyl-2-(2-thienyl)ethyl-4-	
piperidinyl]-N-phenylpropanamide)	
(12) Benzethidine	
(13) Betacetylmethadol	
(14) beta-Hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-	
N-phenylpropanamide)	
(15) beta-Hydroxy-3-methylfentanyl (N-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide)	
(16) beta-Hydroxythiofentanyl (N-[1-[2-hydroxy-2-(thiophen-2-	
yl)ethyl]piperidin-4-yl]-N-phenylpropionamide)	
(17) Betameprodine	
(18) Betamethadol	
(19) beta-Methyl fentanyl (N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-	
yl)propionamide; also known as β -methyl fentanyl)	
(20) beta'-Phenyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N,3-	
diphenylpropanamide; also known as β'-phenyl fentanyl; 3-phenylpropanoyl fentanyl)	
(21) Betaprodine	•
(22) Butyryl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-	-
phenylbutyramide)	
(23) Clonitazene	
$(24) \ \ Crotonyl \ \ fentanyl \ \ ((E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-yl)-N-(24)$	-
enamide)	
(25) Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide)	-
phenylcyclopentanecarboxamide)	•
(26) Cyclopropyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide)	•
(27) Dextromoramide	
(28) Diampromide	
(29) Diethylthiambutene	
(30) Difenoxin	
(31) Dimenoxadol	
(32) Dimepheptanol	
(33) Dimethylthiambutene	
(34) Dioxaphetyl butyrate	
(35) Dipipanone	
(37) Etonitazene	
(38) Etoxeridine	
(39) Fentanyl carbamate (ethyl (1-phenethylpiperidin-4-	
yl)(phenyl)carbamate)	
(40) 4-Fluoroisobutyryl fentanyl (N-(4-fluorophenyl)-N-(1-	-
phenethylpiperidin-4-yl)isobutyramide; also known as para-	-
fluoroisobutyryl fentanyl)	
(41) 2'-Fluoro ortho-fluorofentanyl (N-(1-(2-fluorophenethyl)piperidin-4-	
yl)-N-(2-fluorophenyl)propionamide; also known as 2'-fluoro 2-	
fluorofentanyl)	
(42) Furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide)	
(43) Furethidine	
(44) Hydroxypethidine	

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(45) Isobutyryl	fentanyl	(N-(1-phe	enethylpi	peridin	-4-yl)- <i>N</i> -
phenylisobutyramide)					
(46) Isotonitazene (Abenzimidazol-1-yl)etha					
(47) Ketobemidone					
(48) Levomoramide					
(49) Levomorannue (49) Levophenacylmorpha					
(50) Methoxyacetyl fenta					
phenylacetamide)		, (1 PII)	o11 y 1 p 1		
(51) 4'-Methyl acetyl fer					
N-phenylacetamide)					
(52) 3-Methylfentanyl	(N-[3-methy])	1-1-(2-phe	nylethyl)	-4-pipe	ridyl]- <i>N</i> -
phenylpropanamide)					
(53) 3-Methylthiofentany					
N-phenylpropanamide)					
(54) Morpheridine					
(55) MPPP $(1-methyl-4-ph)$					
(56) MT–45 (1-cyclohexyl- (57) Noracymethadol					
(58) Norlevorphanol					
(59) Normethadone					
(60) Norpipanone					
(61) Ocfentanil (N-(2-fluo	rophenvl)-2-m	ethoxv-N-	-(1-phene	thylnin	eridin-4-
yl)acetamide)					
(62) ortho-Fluoroacı					
phenethylpiperidin-4-y	l)acrylamide)	<i>u</i> =			
(63) ortho-Fluorobut	yryl fenta	anyl	(N-(2-fluc	rophen	yl)- <i>N</i> -(1-
phenethylpiperidin-4-y	l)butyramide:	also kn	own as	2-fluor	butyryl
fentanyl)					
(64) ortho-Fluorofentany	1 (<i>N</i> -(2-fluorop	henyl)- <i>N</i> -	-(1-phene	thylpip	eridin-4-
yl)propionamide); also					
(65) ortho-Fluoroisobi	ıtyryl fent	tanyl	(N-(2-fluc	rophen	yl)- <i>N</i> -(1-
phenethylpiperidin-4-y					
(66) ortho-Methyl					
phenethylpiperidin-4-y					
acetylfentanyl)					
(67) ortho-Methyl i					
methylphenyl)-N-(1-phe					
2-methyl methoxyacety	yı ientanyl)		(37. (4 1. 3	1.	1) 37 (4
(68) para-Chloroisobu phenethylpiperidin-4-y	tyryl fent	anyı ((N-(4-cnlc	ropnen	yı)-N-(1-
pnenetnyipiperidin-4-y.	1)1800utyramic	1e)	(N7 (A £1		1\ \N (1
(69) para-Fluorobuty phenethylpiperidin-4-y	/ryi ienta	ınyı	(1V-(4-IIUC	oropnen	yı)-N-(1-
pneneunyipiperidin-4-y.	1)butyramide)	oronhon		nhonzzi	o+h ₁₇ 1) 4
(70) para-Fluorofentan piperidinyl]propanamic	.yı (1V-(4-11U)	огорпепу.	1)-11-[2-	рпепуг	euny1)-4-
(71) para-Fluoro f					
phenethylpiperidin-4-y					
(72) para-Methoxybut					
phenethylpiperidin-4-y					
(73) para-Methylfentanyl	$(N_{-}(4_{-}methylr))$	nhenvil)-M	(1_nhene	thylnin	eridin_1
yl)propionamide; also l					
(74) PEPAP (1-(2-phenyle					
(75) Phenadoxone					
(76) Phenampromide					
(77) Phenomorphan					
(78) Phenoperidine					
(79) Phenyl fentanyl (N-					
also known as benzoyl					
(80) Piritramide					

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(81) Proheptazine		9643
(82) Properidine		9644
(83) Propiram		9649
(84) Racemoramide		
(85) Tetrahydrofuranyl fentanyl	(N-(1-phenethylpiperidin-4-yl)-N-	0040
phenyltetrahydrofuran-2-carboxamide	Winhons N. M. (9 thions) other 1.4	9843
(86) Thiofentanyl (piperidinyl]propanamide)	(N-phenyl-N-[1-(2-thienyl)ethyl-4-	9835
(87) Thiofuranyl fentanyl phenylthiophene-2-carboxamide; also	(N-(1-phenethylpiperidin-4-yl)-N-	
thiophene fentanyl)		9839
(88) Tilidine		
(89) Trimeperidine		
	2-(dimethylamino)cyclohexyl]- <i>N</i> -	
methylbenzamide)	(N-(1-phenethylpiperidin-4-yl)-N-(
phenylpentanamide)		
(92) Zipeprol (1-methoxy-3-[4-(2-methoxy phenylpropan-2-ol)		
phenyipropan-2-01)	•••••	9013
(c) Opium derivatives. Unless specifi-	contains any of its salts, iso	
cally excepted or unless listed in an-	salts of isomers whenever the	
other schedule, any of the following opium derivatives, its salts, isomers,	of such salts, isomers, and sa	
and salts of isomers whenever the ex-	mers is possible within the chemical designation (for pu	
istence of such salts, isomers, and salts	this paragraph only, the term	
of isomers is possible within the spe-	includes the optical, position	
cific chemical designation:	metric isomers):	ana goo
(1) Acetorphine 9319		7249
(2) Acetyldihydrocodeine 9051	Como trado or other name	
(3) Benzylmorphine	atruntamina: Managa:	
(4) Codeine methylbromide 9070 (5) Codeine-N-Oxide 9053	othyl 1H indolo 9	-
(6) Cyprenorphine	othonomino: 2	2-
(7) Desomorphine	aminobutyl) indole; α-Ε	$\mathrm{T};$
(8) Dihydromorphine 9145	and AET.	
(9) Drotebanol		
(10) Etorphine (except hydro-	phetamine	
chloride salt)		
(11) Heroin	methylphenethylamine;	ι-
(13) Methyldesorphine	0 1	
(14) Methyldihydromorphine 9304		.5-
(15) Morphine methylbromide 9305	dimethoxyphenethylamine.	
(16) Morphine methylsulfonate 9306		
(17) Morphine-N-Oxide	2-(4-bromo-2,5-	
(18) Myrophine 9308 (19) Nicocodeine 9309	difficultory briefly 1)-1-	
(20) Nicomorphine	ammoemane, aipii	
(21) Normorphine	desinethyl DOB, 20-	В,
(22) Pholcodine 9314	NEXUS.	e 7396
(23) Thebacon	Some trade or other name	
(d) Hallucinogenic substances. Unless		io.
specifically excepted or unless listed in		
another schedule, any material, com-	2,5-DMA	
pound, mixture, or preparation, which	(5) 2.5-dimethoxy-	-4-
contains any quantity of the following	ethylamphet-amine	
hallucinogenic substances, or which		

Some trade and other

§ 1308.11 Some trade or other names: DOET (6) 2.5-dimethoxy-4-(n)propylthiophenethylamine (other name: 2C-T-7) (7) 4-methoxyamphetamine Some trade or other names: 4-methoxy- α methylphenethylamine; paramethoxyamphetamine, PMA (8) 5-methoxy-3,4methylenedioxy-amphetamine (9) 4-methyl-2,5-dimethoxy-amphetamine Some trade and other names: 4-methyl-2,5dimethoxy-αmethylphenethylamine; "DOM"; and "STP" (10) 3,4-methylenedioxy amphetamine (11)methylenedioxymethamphetamine (MDMA) 3,4-methylenedioxy-N-(12)ethylamphetamine (also known as N-ethyl-alphamethyl-3,4(methylenedioxy)-N-ethyl phenethylamine, MDA, MDE, MDEA (13)N-hydroxy-3,4methylenedioxyamphetamine (also known as N-hydroxyalpha-methyl-3.4(methylenedioxy)phenethylamine, and N-hydroxy MDA (14) 3,4,5-trimethoxy amphetamine 5-methoxy-N,Ndimethyltryptamine Some trade or other names: 5methoxy-3-[2-(dimethylamino)ethyl]indole; 5-MeO-DMT

Alpha-methyltryptamine (other name: AMT) (17) Bufotenine

(16)

	names: $3-(\beta-$	
	Dimethylaminoethyl)-5-	
	hydroxyindole; 3-(2-	
7348	dimethylaminoethyl)-5-	
7411	indolol; N, N-	
7411	dimethylserotonin; 5-hy-	
	droxy-N,N-	
	dimethyltryptamine;	
	mappine	7404
	(18) Diethyltryptamine	7434
	Some trade and other	
	names: N,N-	
7401	Diethyltryptamine; DET	
	(19) Dimethyltryptamine	7435
7395	Some trade or other names:	
1000	DMT	
	(20) 5-methoxy-N,N-	
	diisopropyltryptamine (other	
	name: 5-MeO-DIPT)	7439
	(21) Ibogaine	7260
	Some trade and other	
	names: 7-Ethyl-	
7400	$6,6\beta,7,8,9,10,12,13$ -	
	octahydro-2-methoxy-6,9-	
7405	methano-5H-pyrido [1',	
1100	2':1,2] azepino [5,4-b]	
	indole; Tabernanthe iboga	
	(22) Lysergic acid diethylamide	7315
	(23) Marihuana	7360
	(24) Mescaline	7381
	(25) Parahexyl—7374; some	
7404	trade or other names: 3-	
	Hexyl-1-hydroxy-7,8,9,10-	
	tetrahydro-6,6,9-trimethyl-	
	6H-dibenzo[b,d]pyran;	
	Synhexyl.	
	(26) Peyote	7415
	Meaning all parts of the	
7402	plant presently classified	
	botanically as Lophophora	
7390	williamsii Lemaire, wheth-	
1000	er growing or not, the	
	seeds thereof, any extract	
	from any part of such	
	plant, and every com-	
	pound, manufacture,	
7431	salts, derivative, mixture,	
	or preparation of such	
7432	plant, its seeds or ex-	
7433	tracts	
	(Interprets 21 USC 812(c),	
	Schedule I(c) (12))	
	(27) N-ethyl-3-piperidyl	
	benzilate	7482
	(28) N-methyl-3-piperidyl	
	benzilate	7484
	(29) Psilocybin	7437
	(30) Psilocyn	7438
	(22) = 22200322	. 250
	3.4	

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	Some trade or other names: 1-(1-phenylcyclohexyl)-	7370	(31) Tetrahydrocannabinols (i) Meaning	(3) (i)
	pyrrolidine, PCPy, PHP		tetrahydrocannabinols, ex-	
	(34) Thiophene analog of		cept as in paragraph (d)(31)(ii)	
	phencyclidine		of this section, naturally con-	
	Some trade or other names:		tained in a plant of the genus	
	1-[1-(2-thienyl)-		Cannabis (cannabis plant), as	
	cyclohexyl]-piperidine, 2-		well as synthetic equivalents of the substances contained	
	thienylanalog of phencyclidine, TPCP,		in the cannabis plant, or in	
	phencyclidine, TPCP, TCP		the resinous extractives of	
	(35) 1-[1-(2-		such plant, and/or synthetic	
	thieny-		substances, derivatives, and	
7473	l)cyclohexyl]pyrrolidine		their isomers with similar	
. = . •	Some other names: TCPy		chemical structure and phar-	
;	(36) 4-methylmethcathinone		macological activity to those	
	(Mephedrone)		substances contained in the	
	(37) 3,4-		plant, such as the following:.]
	methylenedioxypyrovalerone		1 cis or trans	
7535	(MDPV)		tetrahydrocannabinol,	
	(38) 2-(2,5-Dimethoxy-4-		and their optical isomers	
	ethylphenyl)ethanamine (2C-		6 cis or trans	
	E)		tetrahydrocannabinol,	
	(39) 2-(2,5-Dimethoxy-4-		and their optical isomers	
	methylphenyl)ethanamine		3, 4 cis or trans	
	(2C–D)		tetrahydrocannabinol, and its optical isomers	
	(40) 2-(4-Chloro-2,5-			
	dimethoxyphenyl)ethanamine		(Since nomenclature of	
	(2C–C)		these substances is not	
	(41) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine		internationally standard- ized, compounds of these	
	(2C-I)		structures, regardless of	
	(42) 2-[4-(Ethylthio)-2,5-		numerical designation of	
	dimethoxyphenyl]ethanamine		atomic positions cov-	
	(2C-T-2)		ered.)	
	(43) 2-[4-(Isopropylthio)-2,5-		(ii) Tetrahydrocannabinols	
	dimethoxyphenyl]ethanamine		does not include any ma-	
7532	(2C-T-4)		terial, compound, mix-	
	(44) 2-(2,5-		ture, or preparation that	
	Dimethoxypheny-		falls within the definition	
	l)ethanamine (2C–H)		of hemp set forth in 7	
	(45) 2-(2,5-Dimethoxy-4-nitro-		U.S.C. 1639o.	
	phenyl)ethanamine (2C–N)		(32) Ethylamine analog of	
	(46) $2-(2,5-Dimethoxy-4-(n)-1)$	7455	phencyclidine]
7594	propylphenyl)ethanamine		Some trade or other names:	
	(2C-P)		N-ethyl-1- phenylcyclohexylamine,	
	(47) 3,4-Methylenedioxy-N-methylcathinone (Methylone)		(1-	
	(48) (1-pentyl-1 <i>H</i> -indol-3-		phenylcyclohexy-	
	yl)(2,2,3,3-		l)ethylamine, N-(1-	
	tetramethylcyclopropy-		phenylcyclohexy-	
(7144)	l)methanone (UR-144)		l)ethylamine,	
	(49) [1-(5-fluoro-pentyl)-1 <i>H</i> -		cyclohexamine, PCE	
	indol-3-yl](2,2,3,3-		(33) Pyrrolidine analog of	
	tetramethylcyclopropy-	7458	phencyclidine	J
	l)methanone (5-fluoro-UR-144,			
(7011)	XLR11)			

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(50) N-(1-adamantyl)-1-pentyl-		(63) 2-(methylamino)-1-	
1 <i>H</i> -indazole-3-carboxamide (APINACA, AKB48)	(7048)	phenylpentan-1-one (pentedrone)	(1246)
(51) quinolin-8-yl 1-pentyl-1 <i>H</i> -	()	(64) 1-(1,3-benzodioxol-5-yl)-2-	(===+)
indole-3-carboxylate (PB-22;		(methylamino)pentan-1-one	
QUPIC)	(7222)	(pentylone, bk-MBDP)	(7542)
(52) quinolin-8-yl 1-(5-		(65) 4-fluoro-N-methylcathinone	
fluoropentyl)-1 <i>H</i> -indole-3-		(4-FMC; flephedrone)	(1238)
carboxylate (5-fluoro-PB-22;	(7005)	(66) 3-fluoro- <i>N</i> -methylcathinone	(1000)
5F-PB-22)	(7225)	(3-FMC)	(1233)
(53) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-		(67) 1-(naphthalen-2-yl)-2-	
fluorobenzyl)-1 <i>H</i> -indazole-3-		(pyrrolidin-1-yl)pentan-1-one (naphyrone)	(1258)
carboxamide (AB-		(68) <i>alpha-</i>	(1200)
FUBINACA)	(7012)	pyrrolidinobutiophenone (α-	
(54) N-(1-amino-3,3-dimethyl-1-		PBP)	(7546)
oxobutan-2-yl)-1-pentyl-1 <i>H</i> -		(69) N - $(1-amino-3-methyl-1-$	` ′
indazole-3-carboxamide		oxobutan-2-yl)-1-	
(ADB-PINACA)	(7035)	(cyclohexylmethyl)-1 <i>H</i> -inda-	
(55) 2-(4-iodo-2,5-		zole-3-carboxamide (AB-	
dimethoxyphenyl)-N-(2-me- thoxybenzyl)ethanamine (25I-		CHMINACA)	(7031)
NBOMe, 2C-I-NBOMe)	(7538)	(70) N - $(1-amino-3-methyl-1-amino-3$	
(56) 2-(4-chloro-2,5-	(1000)	oxobutan-2-yl)-1-pentyl-1 <i>H</i> -indazole-3-carboxamide (AB-	
dimethoxyphenyl)-N-(2-me-		PINACA)	(7023)
thoxybenzyl)ethanamine		(71) [1-(5-fluoropentyl)-1 <i>H</i> -	(1020)
(25C-NBOMe, 2C-C-NBOMe)	(7537)	indazol-3-yl](naphthalen-1-	
(57) 2-(4-bromo-2,5-		yl)methanone (THJ-2201)	(7024)
dimethoxyphenyl)-N-(2-me-		(72) N-(1-amino-3,3-dimethyl-1-	
thoxybenzyl)ethanamine	(EE00)	oxobutan-2-yl)-1-	
(25B-NBOMe, 2C-B-NBOMe)	(7536)	(cyclohexylmethyl)-1 <i>H</i> -inda-	
(58) Marihuana Extract	7350	zole-3-carboxamide (MAB-	
Meaning an extract con-		CHMINACA; ADB-	(7020)
taining one or more		CHMINACA)	(7032)
cannabinoids that has been derived from any		(73) methyl 2-(1-(5-fluoropentyl)-1 <i>H</i> -indazole-3-	
plant of the genus Can-		carboxamido)-3,3-	
nabis, containing greater		dimethylbutanoate (Other	
than 0.3% delta-9-		names: 5F-ADB; 5F-MDMB-	
tetrahydrocannabinol on		PINACA)	7034
a dry weight basis, other		(74) methyl 2-(1-(5-	
than the separated resin		fluoropentyl)-1 <i>H</i> -indazole-3-	
(whether crude or puri-		carboxamido)-3-	
fied) obtained from the plant.		methylbutanoate (Other names: 5F-AMB)	7033
-		(75) N-(adamantan-1-yl)-1-(5-	1000
(59) 4-methyl-N-ethylcathinone (4-MEC)	(1249)	fluoropentyl)-1 <i>H</i> -indazole-3-	
(60) 4-methyl-alpha-	(1210)	carboxamide (Other names:	
pyrrolidinopropiophenone (4-		5F-APINACA, 5F-AKB48)	7049
MePPP)	(7498)	(76) N - $(1-amino-3,3-dimethyl-1-$	
(61) <i>alpha-</i>	/	oxobutan-2-yl)-1-(4-	
pyrrolidinopentiophenone (α-		fluorobenzyl)-1 <i>H</i> -indazole-3-	
PVP)	(7545)	carboxamide (Other names: ADB-FUBINACA)	7010
(62) 1-(1,3-benzodioxol-5-yl)-2-		ADD-F UDINAUA)	1010
(methylamino)butan-1-one	(DE 44 \		
(butylone, bk-MBDB)	(7541)		

	(88) $1-(4-methoxyphenyl)-N-$		77) methyl 2-(1-	7
	methylpropan-2-amine (other		(cyclohexylmethyl)-1 <i>H</i> -	
	names: para-		indole-3-carboxamido)-3,3-	
(1245)	methoxymethamphetamine, PMMA)		dimethylbutanoate (Other	
(1240)	(89) ethyl 2-(1-(5-fluoropentyl)-	7042	names: MDMB-CHMICA,	
	1 <i>H</i> -indazole-3-carboxamido)-	1042	MMB-CHMINACA)	
	3,3-dimethylbutanoate (other		(8) methyl 2-(1-(4-fluorobenzyl)-1 <i>H</i> -indazole-3-	
7036	name: 5F-EDMB-PINACA)		carboxamido)-3,3-	
	(90) methyl 2-(1-(5-		dimethylbutanoate (Other	
	fluoropentyl)-1 <i>H</i> -indole-3-	7020	names: MDMB-FUBINACA)	
	carboxamido)-3,3-	1020	$\begin{array}{ccc} \text{mattes: WDWB-F CBHVHCH} & \dots \\ \text{(9)} & \text{methyl} & 2-(1-(4-)) \end{array}$	
	dimethylbutanoate (other		fluorobenzyl)-1 <i>H</i> -indazole-3-	
	names: 5F-MDMB-PICA; 5F-		carboxamido)-3-	
7041	MDMB-2201)		methylbutanoate, (FUB-	
	$(91) \qquad N-(adamantan-1-yl)-1-(4-yl)$		AMB, MMB-FUBINACA,	
	fluorobenzyl)-1 <i>H</i> -indazole-3-	(7021)	AMB-FUBINACA)	
	carboxamide (other names:		(0) 1-(1,3-benzodioxol-5-yl)-2-	
	FUB-AKB48; FUB-APINACA;		(ethylamino)propan-1-one	
E0.4E	AKB48 N-(4-	7547	(ethylone)	
7047	FLUOROBENZYL))		31) Naphthalen-1-yl 1-(5-	8
	(92) 1-(5-fluoropentyl)- <i>N</i> -(2-phenylpropan-2-yl)-1 <i>H</i> -inda-		fluoropentyl)-1 <i>H</i> -indole-3-	
	zole-3-carboxamide (other		carboxylate (Other names:	
	names: 5F-CUMYL-PINACA;	7221	NM2201; CBL2201)	
7083	SGT-25)		N-(1-amino-3-methyl-1-	
1000	(93) (1-(4-fluorobenzyl)-1 <i>H</i> -		oxobutan-2-yl)-1-(5-	
	indol-3-yl)(2,2,3,3-		fluoropentyl)-1 <i>H</i> -indazole-3-	
	tetramethylcyclopropy-	5 005	carboxamide (Other name:	
	l)methanone (other name:	7025	5F-AB-PINACA)	
7014	FUB-144)		3) 1-(4-cyanobutyl)- <i>N</i> -(2-	
	(94) N-Ethylhexedrone (Other		phenylpropan-2-yl)-1 <i>H</i> -inda-	
	names: α -		zole-3-carboxamide (Other	
	ethylaminohexanophenone; 2-		names: 4-CN-CUMYL-	
	(ethylamino)-1-phenylhexan-		BUTINACA; 4-cyano-CUMYL-BUTINACA; 4-CN-CUMYL	
7246	1-one)		BINACA; CUMYL-4CN-	
	(95) <i>alpha-</i>	7089	BINACA; SGT-78)	
	Pyrrolidinohexanophenone	1000	(4) methyl 2-(1-	
	(Other names: α -PHP; α -		(cyclohexylmethyl)-1 <i>H</i> -	
	pyrrolidinohexanophenone; 1-		indole-3-carboxamido)-3-	
7544	phenyl-2-(pyrrolidin-1- yl)hexan-1-one)		methylbutanoate (Other	
1011	(96) 4 -Methyl- $alpha$ -		names: MMB-CHMICA; AMB-	
	ethylaminopentiophenone	7044	CHMICA)	
	(Other names: 4-MEAP; 2-		1-(5-fluoropentyl)-N-(2-	8
	(ethylamino)-1-(4-		phenylpropan-2-yl)-1 <i>H</i> -	
7245	methylphenyl)pentan-1-one)		pyrrolo[2,3-b]pyridine-3-	
	(97) 4'-Methyl-alpha-		carboxamide (Other name:	
	pyrrolidinohexiophenone	7085	5F-CUMYL-P7AICA)	
	(Other names: MPHP; 4'-		66) N-ethylpentylone (Other	
	methyl-alpha-		names: ephylone, 1-(1,3-	
	pyrrolidinohexanophenone; 1-	40	benzodioxol-5-yl)-2-	
	(4-methylphenyl)-2-	7543	(ethylamino)pentan-1-one)	
7446	(pyrrolidin-1-yl)hexan-1-one)		7) methyl 2-(1-(4-fluorobutyl)-	
	(98) alpha-		1 <i>H</i> -indazole-3-carboxamido)-	
	Pyrrolidinoheptaphenone (Other names: PV8; 1-phenyl-		3,3-dimethylbutanoate (4F–	
	2-(pyrrolidin-1-yl)heptan-1-	7043	MDMB-BINACA, 4F-MDMB-BUTINACA)	
7548	one)	GEUI	DOTIMAUA)	

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(99) 4'-Chloro-alpha-	Some trade or other names:
pyrrolidinovalerophenone	2-amino-1-phenyl-1-
(Other names: 4-chloro-α-	propanone, alpha-
PVP; 4'-chloro-α-	aminopropiophenone, 2-
pyrrolidinopentiophenone; 1-	aminopropiophenone, and
(4-chlorophenyl)-2- (pyrrolidin-1-yl)pentan-1-one) 7443	norephedrone
(100) 2-(ethylamino)-2-(3-	(5) 4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4-methyl-
methoxyphenyl)cyclohexan-1-	5-(4-methylphenyl)-2-
one (methoxetamine, MXE) 7286	oxazolamine; 4-methyl-5-(4-
·	methylphenyl)-4,5-dihydro-
(e) Depressants. Unless specifically	1,3-oxazol-2-amine) 1595
excepted or unless listed in another	(6) Fenethylline 1503
schedule, any material, compound,	(7) Mesocarb (<i>N</i> -phenyl- <i>N'</i> -(3-(1-
mixture, or preparation which contains	phenylpropan-2-yl)-1,2,3-
any quantity of the following sub-	oxadiazol-3-ium-5-
stances having a depressant effect on the central nervous system, including	yl)carbamimidate) 1227
its salts, isomers, and salts of isomers	(8) Methcathinone (Some other
whenever the existence of such salts,	names: 2-(methylamino)-
isomers, and salts of isomers is possible	propiophenone; alpha-
within the specific chemical designa-	(methylamin- o)propiophenone: 2-
tion:	o)propiophenone; 2- (methylamino)-1-
(1) gamma-hydroxybutyric acid	phenylpropan-1-one; alpha-N-
(some other names include	methylaminopropiophenone;
GHB; gamma-	monomethylpropion;
hydroxybutyrate; 4-	ephedrone; N-
hydroxybutyrate; 4-	methylcathinone;
hydroxybutanoic acid; so-	methylcathinone; AL-464;
dium oxybate; sodium	AL-422; AL-463 and UR1432),
oxybutyrate) 2010	its salts, optical isomers and
(2) Mecloqualone 2572	salts of optical isomers 1237
(3) Methaqualone 2565	(9) Methiopropamine (<i>N</i> -methyl-1-(thiophen-2-
(f) Stimulants. Unless specifically ex-	yl)propan-2-amine) 1478
cepted or unless listed in another	(10) (\pm) cis -4-methylaminorex
schedule, any material, compound,	$(\pm)cis$ -1-incony minimized $((\pm)cis$ -4,5-dihydro-4-methyl-5-
mixture, or preparation which contains	phenyl-2-oxazolamine) 1590
any quantity of the following sub-	(11) N-ethylamphetamine 1475
stances having a stimulant effect on	(12) N,N-dimethylamphetamine
the central nervous system, including	(also known as N,N-alpha-
its salts, isomers, and salts of isomers:	trimethyl-
(1) Amineptine (7-[(10,11-	benzeneethanamine; N,N -
dihydro-5H-	alpha-
dibenzo[a,d]cyclohepten-5-	trimethylphenethylamine) 1480
yl)amino]heptanoic acid) 1219	(g) Cannabimimetic agents. Unless spe-
(2) Aminorex (Some other	cifically exempted or unless listed in
names: aminoxaphen; 2-	another schedule, any material, com-
amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenly-2-	pound, mixture, or preparation which
oxazolamine) 1585	contains any quantity of the following
(3) N-Benzylpiperazine (some	substances, or which contains their
other names: BZP, 1-	salts, isomers, and salts of isomers
benzylpiperazine) 7493	whenever the existence of such salts,
(4) Cathinone	isomers, and salts of isomers is possible
	within the specific chemical designa-
	tion:

Drug Enforcement Administration, Justice

(1) 5-(1,1-dimethylheptyl)-2- [(1R,3S)-3- hydroxycyclohexyl]-phenol (CP-47,497)	297	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: (A) Replacement of the phenyl por-
(3) 1-pentyl-3-(1-naph-thoyl)indole (JWH-018 and	298	tion 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,
AM678)	.18	alkoxyl, hydroxyl, halo, haloalkyl, amino or nitro groups;
(5) 1-hexyl-3-(1-naph-	.73	(C) Substitution in or on the piperidine ring with alkyl, alkenyl, alkoxyl,
(6) 1-[2-(4-morpholinyl)ethyl]-3-)19	ester, ether, hydroxyl, halo, haloalkyl, amino or nitro groups;
(7) 1-pentyl-3-(2-	200	(D) Replacement of the aniline ring with any aromatic monocycle whether or not further substituted in or on the
methoxyphenylacetyl)indole (JWH-250)	250	aromatic monocycle; and/or (E) Replacement of the N -propionyl
methoxynaphthoyl)]indole (JWH-081)70	081	group by another acyl group. (ii) This definition includes, but is not limited to, the following sub-
(9) 1-pentyl-3-(4-methyl-1-naph- thoyl)indole (JWH-122)	.22	stances: (A)–(B) [Reserved]
	898	(31)-(49) [Reserved]. (50) 2-(2-(4-butoxybenzyl)-5- nitro-1 <i>H</i> -benzimidazol-1-yl)-
	201	N,N-diethylethan-1-amine, its isomers, esters, ethers, salts,
	94	and salts of isomers, esters and ethers (Other name:
(14) 1-cyclohexylethyl-3-(2-	.04	$\begin{array}{cccccccccccccccccccccccccccccccccccc$
	800	diethylethan-1-amine, its isomers, esters, ethers, salts,
(15) 1-pentyl-3-(2- chlorophenylacetyl)indole (JWH-203)	203	and salts of isomers, esters and ethers (Other names: Etodesnitazene; etazene) 9765
(h) Temporary listing of substances su ject to emergency scheduling. Any mat rial, compound, mixture or preparatic which contains any quantity of the following substances:	e- on	(52) N,N-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of iso-
(1)–(29) [Reserved]. (30) Fentanyl-related substances, their isomers, esters, ethers, salts and salts of isomers, esters and ethers	350	mers, esters and ethers (Other name: Flunitazene) 9756 (53) N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters,
(i) Fentanyl-related substance mea any substance not otherwise list under another Administration Co	ed n-	ethers, salts, and salts of isomers, esters and ethers (Other name:
trolled Substance Code Number, as for which no exemption or approval		Metodesnitazene) 9764

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(54) N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (Other name: Metonitazene) ..
(55) 2-(4-ethoxybenzyl)-5-nitro-

(55) 2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1*H*-benzimidazole, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (Other names: *N*pyrrolidino etonitazene; etonitazepyne)

(56) N,N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine, its isomers, esters, ethers, salts, and salts of isomers, esters and ethers (Other name: Protonitazene)

[39 FR 22141, June 20, 1974]

EDITORIAL NOTE: For FEDERAL REGISTER citations affecting §1308.11, see the List of CFR Sections Affected, which appears in the Finding Aids section of the printed volume and at www.govinfo.gov.

EFFECTIVE DATE NOTES:

1.At 83 FR 5191, Feb. 6, 2018, \$1308.11 was amended by adding paragraph (h)(30), effective Feb. 6, 2018, through Feb. 6, 2020. Effective Feb. 6, 2020, Congress extended the effective period for paragraph (h)(30) until May 6,

2021, by Public Law 116-114, Effective May 4. 2021, Congress extended the effective period for paragraph (h)(30) until October 22, 2021, by Public Law 117-12. Effective Sept. 30, 2021, Congress extended the effective period for paragraph (h)(30) until Jan. 28, 2022, by Public Law 117-43. Effective Jan. 13, 2022. Congress extended the effective period for paragraph (h)(30) until Feb. 18, 2022, by Public Law 117-70. Effective Feb. 18, 2022, Congress extended the effective period for paragraph (h)(30) until Mar. 11, 2022, by Public Law 117-86. Effective Mar. 11, 2022, Congress extended the effective period for paragraph (h)(30) until Mar. 15, 2022 by Public Law 117-95. Effective Mar. 15, 2022, Congress extended the effective period for paragraph (h)(30) until Dec. 31, 2022 by Public Law No. 117-103. Effective Dec. 29, 2022, Congress extended the effective period for paragraph (h)(30) until Dec. 31, 2024 by Public Law No. 117-328.

2. At 87 FR 21561, Apr. 12, 2022, §1308.11 was amended by adding paragraphs (h)(50) through (h)(56), effective Apr. 12, 2022 through Apr. 12, 2024.

3. At 88 FR 13694, Mar. 6, 2023, §1308.11 was amended by redesignating paragraphs (b)(22) through (93) as paragraphs (b)(23) through (94), respectively and adding new paragraph (b)(22) and removing and reserving paragraph (h)(49), effective Apr. 5, 2023. For the convenience of the user, the added text is set forth as follows:

§ 1308.11 Schedule I.

* * * * * * * * *

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§ 1308.12 Schedule II.

(a) 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.

(b) Substances, vegetable origin or chemical synthesis. Unless specifically excepted or unless listed in another schedule, 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:

(1) Opium and opiate, and any salt, compound, derivative, or preparation of opium or opiate excluding apomorphine, thebaine-derived butorphanol, dextrorphan, nalbuphine, naldemedine, nalmefene, naloxegol, naloxone, 6β-naltrexol, naltrexone, and samidorphan, and their respective salts, but including the following:

9098

(i) Codeine	9050
(ii) Dihydroetorphine	9334
(iii) Ethylmorphine	9190
(iv) Etorphine hydrochloride	9059
(v) Granulated opium	9640
(vi) Hydrocodone	9193
(vii) Hydromorphone	9150
(viii) Metopon	9260
(ix) Morphine	9300
(x) Noroxymorphone	9668

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