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§ 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 to §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 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 paragraph (b)(34) only, the term isomer includes the optical and geometric isomers):

(1) Acetyl-alpha-methylfentanyl	
(N-[1-(1-methyl-2-phenethyl)-4-	
piperidinyl]- <i>N</i> -	
phenylacetamide)	9815
(2) Acetylmethadol	9601
(3) Allylprodine	9602
(4) Alphacetylmethadol (except	
levo-alphacetylmethadol also	
known as levo-alpha-	
acetylmethadol, levomethadyl	
acetate, or LAAM)	9603
(5) Alphameprodine	9604
(6) Alphamethadol	9605
(7) Alpha-methylfentanyl (N-[1-	
(alpha-methyl-beta-	
phenyl)ethyl-4-piperidyl]	
propionanilide; 1-(1-methyl-2-	
phenylethyl)-4-(N-propanilido)	
piperidine)	9814
(8) Alpha-methylthiofentanyl	
(N-[1-methyl-2-(2-	
thienyl)ethyl-4-piperidinyl]- <i>N</i> -	
phenylpropanamide)	9832
(9) Benzethidine	9606
(10) Betacetylmethadol	9607
(11) Beta-hydroxyfentanyl (N-[1-	
(2-hydroxy-2-phenethyl)-4-	
piperidinyl]-N-	0000
phenylpropanamide)	9830
(12) Beta-hydroxy-3-	
methylfentanyl (other name:	
N-[1-(2-hydroxy-2-phenethyl)-	
3-methyl-4-piperidinyl]-N-	0001
phenylpropanamide	9831
(13) Betameprodine	9608
(14) Betamethadol	9609
(15) Betaprodine	9611
(16) Clonitazene	9612
(17) Dextromoramide	9613
(18) Diampromide	9615
(19) Diethylthiambutene	9616
(20) Difference del	9168
(21) Dimenoxadol	9617 9618
(22) Dimepheptanol	9619
(24) Dieverhetyl bytyrote	9621
(24) Dioxaphetyl butyrate	
(25) Dipipanone	9622
(26) Ethylmethylthiambutene	9623 9624
(27) Etonitazene	
(28) Etoxeridine	9625 9626
(29) Furethidine	9626

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(30) Hydroxypethidine	9627	(11) Heroin 92	200
(31) Ketobemidone	9628		301
(32) Levomoramide	9629		302
(33) Levophenacylmorphan	9631	· · · · · ·	304
(34) 3-Methylfentanyl $(N-[3-$			305
methyl-1-(2-phenylethyl)-4-		· · · · -	306
piperidyl]-N-	0010	· · · · · · · · · · · · · · · · · · ·	307
phenylpropanamide)	9813	· · · · · · =	308
methyl-1-(2-thienyl)ethyl-4-			309
piperidinyl]-N-		(=-, =-=	312
phenylpropanamide)	9833	-	313
(36) Morpheridine	9632	-	314
(37) MPPP (1-methyl-4-phenyl-4-		, , , , , , , , , , , , , , , , , , , ,	315
propionoxypiperidine)	9661		
(38) Noracymethadol	9633	(d) Hallucinogenic substances. Unle	
(39) Norlevorphanol	9634	specifically excepted or unless listed	
(40) Normethadone	9635	another schedule, any material, co	
(41) Norpipanone	9636	pound, mixture, or preparation, whi	
(42) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-		contains any quantity of the followi hallucinogenic substances, or whi	-
phenethyl)-4-piperidinyl]		contains any of its salts, isomers, a	
propanamide	9812	salts of isomers whenever the existen	
(43) PEPAP (1-(-2-phenethyl)-4-	0012	of such salts, isomers, and salts of is	
phenyl-4-acetoxypiperidine	9663	mers is possible within the speci	
(44) Phenadoxone	9637	chemical designation (for purposes	
(45) Phenampromide	9638	this paragraph only, the term "isome	
(46) Phenomorphan	9647	includes the optical, position and ge	
(47) Phenoperidine	9641	metric isomers):	
(48) Piritramide	9642 9643	,	
(49) Proheptazine(50) Properidine	9643 9644	(1) Alpha-ethyltryptamine 72	249
(51) Propiram	9649	Some trade or other names:	
(52) Racemoramide	9645	etryptamine; Monase; α -	
(53) Thiofentanyl (N-phenyl-N-[1-		ethyl-1H-indole-3-	
(2-thienyl)ethyl-4-piperidinyl]-		ethanamine; 3-(2-	
propanamide	9835	aminobutyl) indole; α -ET;	
(54) Tilidine	9750	and AET.	
(55) Trimeperidine	9646	(2) 4-bromo-2,5-dimethoxy-am-	201
(c) Opium derivatives. Unless spe	ecifi-	2	391
cally excepted or unless listed in	an-	Some trade or other names:	
other schedule, any of the follo	wing	4-bromo-2,5-dimethoxy-α- methylphenethylamine; 4-	
opium derivatives, its salts, isor		bromo-2,5-DMA	
and salts of isomers whenever the		(3) 4-Bromo-2,5-	
istence of such salts, isomers, and			392
of isomers is possible within the cific chemical designation:	spe-	Some trade or other names:	
cific chemical designation.		2-(4-bromo-2,5-	
(1) Acetorphine	9319	dimethoxyphenyl)-1-	
(2) Acetyldihydrocodeine	9051	aminoethane; alpha-	
(3) Benzylmorphine	9052	desmethyl DOB; 2C-B,	
(4) Codeine methylbromide	9070	Nexus.	
(5) Codeine-N-Oxide	9053	. , ,	396
(6) Cyprenorphine	9054	Some trade or other names:	
(7) Desomorphine	9055	2,5-dimethoxy-α-	
(8) Dihydromorphine	9145	methylphenethylamine;	
(9) Drotebanol	9335	2,5-DMA	
(10) Etorphine (except hydro- chloride salt)	0056	(5) 2,5-dimethoxy-4-	399
omorine sam,	9056	ethylamphet-amine	บฮฮ

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Some trade or other names: DOET		Some trade and other names: 3-(β-	
(6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine		Dimethylaminoethyl)-5- hydroxyindole; 3-(2-	
(other name: 2C-T-7)	7348	dimethylaminoethyl)-5-	
(7) 4-methoxyamphetamine	7411	indolol; N, N-	
Some trade or other names:		dimethylserotonin; 5-hy- droxy-N,N-	
4-methoxy-α-		dimethyltryptamine;	
methylphenethylamine; paramethoxyamphetamin-		mappine	
e, PMA		(18) Diethyltryptamine	7434
(8) 5-methoxy-3,4-		Some trade and other	
methylenedioxy-amphetamine	7401	names: N,N-	
(9) 4-methyl-2,5-dimethoxy-am-		Diethyltryptamine; DET	
phetamine	7395	(19) Dimethyltryptamine	7435
Some trade and other names: 4-methyl-2,5-		Some trade or other names: DMT	
dimethoxy-α-		(20) 5-methoxy-N,N-	
methylphenethylamine;		diisopropyltryptamine (other name: 5-MeO-DIPT)	7439
"DOM"; and "STP" (10) 3,4-methylenedioxy amphet-		(21) Ibogaine	7260
amine	7400	Some trade and other	1200
(11) 3,4-	1 100	names: 7-Ethyl-	
methylenedioxymethampheta-		6,6β,7,8,9,10,12,13-	
mine (MDMA)	7405	octahydro-2-methoxy-6,9-	
(12) 3,4-methylenedioxy-N-		methano-5H-pyrido [1',	
ethylamphetamine (also		2':1,2] azepino [5,4-b]	
known as N-ethyl-alpha-meth-		indole; Tabernanthe iboga (22) Lysergic acid diethylamide	7315
yl-3,4(methylenedioxy)- phenethylamine, N-ethyl		(23) Marihuana	7360
MDA, MDE, MDEA	7404	(24) Mescaline	7381
(13) N-hydroxy-3,4-	. 101	(25) Parahexyl—7374; some trade	1001
methylenedioxyamphetamine		or other names: 3-Hexyl-1-hy-	
(also known as N-hydroxy-		droxy-7,8,9,10-tetrahydro-6,6,9-	
alpha-methyl-		trimethyl-6H-	
3,4(methylenedioxy)- phenethylamine, and N-hy-		dibenzo[b,d]pyran; Synhexyl.	P.415
droxy MDA	7402	(26) Peyote	7415
(14) 3,4,5-trimethoxy amphet-	1102	Meaning all parts of the plant presently classified	
amine	7390	botanically as Lophophora	
(15) 5-methoxy-N,N-		williamsii Lemaire, whether	
dimethyltryptamine Some		growing or not, the seeds	
trade or other names: 5-		thereof, any extract from	
methoxy-3-[2- (dimethylamino)ethyl]indole;		any part of such plant, and every compound, manufac-	
5-MeO-DMT	7431	ture, salts, derivative,	
(16) Alpha-methyltryptamine	1101	mixture, or preparation of	
(other name: AMT)	7432	such plant, its seeds or ex-	
(17) Bufotenine	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
		(21) Totrobydroconnobinola	7270

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Meaning		Some other names: TCPy	
tetrahydrocannabinols		(36) 4-methylmethcathinone	
naturally contained in a		, 1	248
plant of the genus Can-		(37) 3,4-	
nabis (cannabis plant), as		methylenedioxypyrovalerone	
well as synthetic equiva-			535
lents of the substances		(38) 2-(2,5-Dimethoxy-4-	
contained in the cannabis		ethylphenyl)ethanamine (2C-	
plant, or in the resinous			509
extractives of such plant,		(39) 2-(2,5-Dimethoxy-4-	
and/or synthetic sub-		methylphenyl)ethanamine	- 00
stances, derivatives, and			508
their isomers with similar		(40) 2-(4-Chloro-2,5-	
chemical structure and		dimethoxyphenyl)ethanamine	-10
pharmacological activity			519
to those substances con-		(41) 2-(4-Iodo-2,5-	
tained in the plant, such		dimethoxyphenyl)ethanamine	-10
as the following:			518
1 cis or trans			
tetrahydrocannabinol, and		dimethoxyphenyl]ethanamine	385
their optical isomers 6 cis or trans		(2C-T-2)	ာဝပ
6 cis or trans tetrahydrocannabinol, and		dimethoxyphenyl]ethanamine	
their optical isomers			532
3, 4 cis or trans		(44) 2-(2,5-	102
tetrahydrocannabinol, and		Dimethoxyphenyl)ethanamine	
its optical isomers			517
(Since nomenclature of		(45) 2-(2,5-Dimethoxy-4-nitro-	, . .
these substances is not		· · · · · · · · · · · · · · · · · · ·	521
internationally standard-		(46) 2-(2,5-Dimethoxy-4-(n)-	
ized, compounds of these		propylphenyl)ethanamine (2C–	
structures, regardless of			524
numerical designation of		(47) 3,4-Methylenedioxy-N-	
atomic positions covered.)			540
(32) Ethylamine analog of			
phencyclidine	7455	(e) Depressants. Unless specifical excepted or unless listed in anoth	
Some trade or other names:		schedule, any material, compour	
N-ethyl-1-		mixture, or preparation which contain	
phenylcyclohexylamine,		any quantity of the following su	
(1-		stances having a depressant effect	
phenylcyclohexy-		the central nervous system, including	
l)ethylamine, N-(1-		its salts, isomers, and salts of isome	_
phenylcyclohexy-		whenever the existence of such salt	
l)ethylamine,		isomers, and salts of isomers is possib	-
cyclohexamine, PCE		within the specific chemical design	
(33) Pyrrolidine analog of		tion:	
phencyclidine	7458		
Some trade or other names:		(1) gamma-hydroxybutyric acid	
1-(1-phenylcyclohexyl)-		(some other names include	
pyrrolidine, PCPy, PHP		GHB; gamma-	
(34) Thiophene analog of	E 4E0	hydroxybutyrate; 4-	
phencyclidine	7470	hydroxybutyrate; 4-	
Some trade or other names:		hydroxybutanoic acid; sodium	
1-[1-(2-thienyl)-			010
cyclohexyl]-piperidine, 2-			572
thienylanalog of phencyclidine, TPCP, TCP		• •	565
(35) phencychaine, TPCP, TCP		(f) Stimulants. Unless specifically e	v.
thienyl)cyclohexyl]pyrrolidine	7473	cepted or unless listed in anoth	
omion's 1/0's oromov's 11h's 11 origine	1110	cobood of annops moned in which	.01

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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:

(1) Aminorex (Some other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenly-2-	
oxazolamine)	1585
(2) N-Benzylpiperazine (some	
other names: BZP, 1-	
benzylpiperazine)	7493
(3) Cathinone	1235
Some trade or other names:	1200
2-amino-1-phenyl-1-	
propanone, alpha-	
aminopropiophenone, 2-	
aminopropiophenone, and	
norephedrone	
(4) Fenethylline	1503
(5) Methcathinone (Some other	1505
names: 2-(methylamino)-	
propiophenone; alpha-	
(methylamino)propiophenone;	
2-(methylamino)-1-	
phenylpropan-1-one; alpha-N-	
methylaminopropiophenone;	
monomethylpropion;	
ephedrone; N-	
methylcathinone;	
methylcathinone; AL-464; AL-	
422; AL–463 and UR1432), its	
salts, optical isomers and salts	1005
of optical isomers	1237
(6) (\pm) cis -4-methylaminorex	
$((\pm)cis$ -4,5-dihydro-4-methyl-5-	
phenyl-2-oxazolamine)	1590
(7) N-ethylamphetamine	1475
(8) N,N -dimethylamphetamine	
(also known as N,N-alpha-	
trimethyl-benzeneethanamine;	
N,N-alpha-	
trimethylphenethylamine)	1480

(g) Cannabimimetic agents. Unless 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:

$(1) \qquad 5-(1,1-dimethylheptyl)-2-$	
[(1R,3S)-3-hydroxycyclohexyl]-	7007
phenol (CP-47,497)	7297
(2) 5-(1,1-dimethyloctyl)-2-	
[(1R,3S)-3-hydroxycyclohexyl]-	
phenol (cannabicyclohexanol	7000
or CP-47,497 C8-homolog) (3) 1-pentyl-3-(1-naph-	7298
thoyl)indole (JWH–018 and	
AM678)	7118
(4) 1-butyl-3-(1-naphthoyl)indole	1110
(JWH-073)	7173
(5) 1-hexyl-3-(1-naphthoyl)indole	1110
(JWH–019)	7019
(6) 1-[2-(4-morpholinyl)ethyl]-3-	1010
(1-naphthoyl)indole (JWH–200)	7200
(7) 1-pentyl-3-(2-	1200
methoxyphenylacetyl)indole	
(JWH–250)	6250
(8) 1-pentyl-3-[1-(4-	
methoxynaphthoyl)]indole	
(JWH-081)	7081
(9) 1-pentyl-3-(4-methyl-1-naph-	
thoyl)indole (JWH-122)	7122
(10) 1-pentyl-3-(4-chloro-1-naph-	
thoyl)indole (JWH-398)	7398
(11) 1-(5-fluoropentyl)-3-(1-naph-	
thoyl)indole (AM2201)	7201
(12) 1-(5-fluoropentyl)-3-(2-	
iodobenzoyl)indole (AM694)	7694
(13) 1-pentyl-3-[(4-methoxy)-ben-	5 101
zoyl]indole (SR-19 and RCS-4)	7104
(14) 1-cyclohexylethyl-3-(2-	
methoxyphenylacetyl)indole	7000
7008 (SR–18 and RCS–8)	7008
(15) 1-pentyl-3-(2-chlorophenylacetyl)indole	
(JWH-203)	7203
· · · · · · · · · · · · · · · · · · ·	
(h) Tomporary listing of substances	euh_

- (h) Temporary listing of substances subject to emergency scheduling. Any material, compound, mixture or preparation which contains any quantity of the following substances:
- $\begin{array}{ccc} (1) & 3,4\text{-methylenedioxy-N-} \\ \text{methylcathinone} & (\text{Other names:} \\ \text{methylone}) 7540 \end{array}$
- (2) 5-(1,1-Dimethyloctyl)-2-[(1*R*,3*S*)-3-hydroxycyclohexyl]-phenol, its optical, positional, and geometric isomers, salts and salts of isomers—7298 (Other names: cannabicyclohexanol and CP-47,497 C8 homologue)
- (3) 1-Butyl-3-(1-naphthoyl)indole, its optical, positional, and geometric isomers, salts and salts of isomers—7173 (Other names: JWH-073)
- (4) 1-[2-(4-Morpholinyl)ethyl]-3-(1-naphthoyl)indole, its optical, positional, and geometric isomers, salts

- and salts of isomers—7200 (Other names: JWH-200)
- (5) 1-Pentyl-3-(1-naphthoyl)indole, its optical, positional, and geometric isomers, salts and salts of isomers—7118 (Other names: JWH-018 and AM678)
- (6) 4-methyl-N-methylcathinone—1248 (Other names: mephedrone)
- (7) 3,4-methylenedioxy-N-methylcathinone—7540 (Other names: methylone)
- (8) 3,4-methylenedioxypyrovalerone—7535 (Other names: MDPV)
- (9) (1-pentyl-1*H*-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone, its optical, positional, and geometric isomers, salts and salts of isomers—7144 (Other names: UR-144, 1-pentyl-3-(2,2,3,3-

tetramethylcyclopropoyl)indole)

- (10) [1-(5-fluoro-pentyl)-1H-indol-3-yl](2,2,3,3-
- tetramethylcyclopropyl)methanone, its optical, positional, and geometric isomers, salts and salts of isomers—7011 (Other names: 5-fluoro-UR-144, 5-F-UR-144, XLR11, 1-(5-fluoro-pentyl)-3-(2,2,3,3-

tetramethylcyclopropoyl)indole)

- (11) N-(1-adamantyl)-1-pentyl-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers—7048 (Other names: APINACA, AKB48)
- (12) 2-(4-iodo-2,5-dimethoxyphenyl)-*N*-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts and salts of isomers—7538 (Other names: 25I–NBOMe; 2C–I–NBOMe; 25I; Cimbi-5)
- (13) 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts and salts of isomers—7537 (Other names: 25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82)
- (14) 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, its optical, positional, and geometric isomers, salts and salts of isomers—7536
- (Other names: 25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36)
- (15) Quinolin-8-yl 1-pentyl-1*H*-indole-3-carboxylate, its optical, positional, and geometric isomers, salts and salts of isomers—7222 (Other names: PB-22; QUPIC)
- (16) Quinolin-8-yl 1-(5-fluoropentyl)-1*H*-indole-3-carboxylate, its optical, po-

- sitional, and geometric isomers, salts and salts of isomers—7225 (Other names: 5-fluoro-PB-22; 5F-PB-22)
- (17) N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers—7012 (Other names: AB-FUBINACA)
- (18) N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1*H*-indazole-3-carboxamide, its optical, positional, and geometric isomers, salts and salts of isomers—7035 (Other names: ADB-PINACA)
- (19) 4-methyl-*N*-ethylcathinone, its optical, positional, and geometric isomers, salts and salts of isomers—1249 (Other names: 4-MEC; 2-(ethylamino)-1-(4-methylphenyl)propan-1-one)
- (20) 4-methyl-alpha-pyrrolidinopropiophenone, its optical, positional, and geometric isomers, salts and salts of isomers—7498 (Other names: 4-MePPP; MePPP; 4-methyl-α-pyrrolidinopropiophenone; 1-(4-methylphenyl)-2-(pyrrolidin-1-yl)-propan-1-one)
- (21) alpha-pyrrolidinopentiophenone, its optical, positional, and geometric isomers, salts and salts of isomers—7545 (Other names: α -PVP; α -pyrrolidinovalerophenone; 1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one)
- (22) Butylone, its optical, positional, and geometric isomers, salts and salts of isomers—7541 (Other names: bk-MBDB; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one)
- (23) Pentedrone, its optical, positional, and geometric isomers, salts and salts of isomers—1246 (Other names: α-methylaminovalerophenone; 2-(methylamino)-1-phenylpentan-1-one)
- (24) Pentylone, its optical, positional, and geometric isomers, salts and salts of isomers—7542 (Other names: bk-MBDP; 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one)
- (25) 4-fluoro-*N*-methylcathinone, its optical, positional, and geometric isomers, salts and salts of isomers—1238 (Other names: 4-FMC; flephedrone; 1-(4-fluorophenyl)-2-(methylamino)propan-1-one)
- (26) 3-fluoro-N-methylcathinone, its optical, positional, and geometric isomers, salts and salts of isomers—1233 (Other names: 3-FMC; 1-(3-

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fluorophenyl)-2-(methylamino)propan-1-one)

(27) Naphyrone, its optical, positional, and geometric isomers, salts and salts of isomers—1258 (Other names: naphthylpyrovalerone; 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one)

(28) alpha-pyrrolidinobutiophenone, its optical, positional, and geometric isomers, salts and salts of isomers—7546 (Other names: α -PBP; 1-phenyl-2-(pyrrolidin-1-yl)butan-1-one)

[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.fdsys.gov.

§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, nalmefene, naloxone, and naltrexone, and their respective salts, but including the following:

(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) Opium extracts	9610
(xi) Opium fluid	9620

(xii) Oripavine	9330
(xiii) Oxycodone	9143
(xiv) Oxymorphone	9652
(xv) Powdered opium	9639
(xvi) Raw opium	9600
(xvii) Thebaine	9333
(xviii) Tincture of opium	9630

- (2) Any salt, compound, derivative, or preparation thereof which is chemically equivalent or identical with any of the substances referred to in paragraph (b) (1) of this section, except that these substances shall not include the isoquinoline alkaloids of opium.
 - (3) Opium poppy and poppy straw.
- (4) 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 equivalent or identical with any of these substances, except that the substances shall not include decocainized coca leaves or extraction of coca leaves, which extractions do not contain cocaine or ecgonine.
- (5) 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.
- (c) 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:

(1) Alfentanil	9737
(2) Alphaprodine	9010
(3) Anileridine	9020
(4) Bezitramide	9800
(5) Bulk dextropropoxyphene	
(non-dosage forms)	9273
(6) Carfentanil	9743
(7) Dihydrocodeine	9120
(8) Diphenoxylate	9170
(9) Fentanyl	9801
(10) Isomethadone	9226
(11) Levo-alphacetylmethadol	9648