SENATE BILL No. 234

DIGEST OF INTRODUCED BILL

Citations Affected: IC 35-41-1-26.3; IC 35-48-2-4.

Synopsis: Synthetic drugs (including "bath salts"). Adds additional chemical compounds (including some compounds sold as "bath salts") to the definition of synthetic cannabinoids, and expands the definition of synthetic cannabinoids to encompass certain chemical compounds that are structurally related to synthetic cannabinoids.

Effective: July 1, 2012.

Alting, Merritt

January 4, 2012, read first time and referred to Committee on Corrections, Criminal, and Civil Matters.

Second Regular Session 117th General Assembly (2012)

PRINTING CODE. Amendments: Whenever an existing statute (or a section of the Indiana Constitution) is being amended, the text of the existing provision will appear in this style type, additions will appear in **this style type**, and deletions will appear in ~~this style type.~~

Additions: Whenever a new statutory provision is being enacted (or a new constitutional provision adopted), the text of the new provision will appear in **this style type**. Also, the word **NEW** will appear in that style type in the introductory clause of each SECTION that adds a new provision to the Indiana Code or the Indiana Constitution.

Conflict reconciliation: Text in a statute in *this style type* or ~~*this style type*~~ reconciles conflicts between statutes enacted by the 2011 Regular Session of the General Assembly.

SENATE BILL No. 234

A BILL FOR AN ACT to amend the Indiana Code concerning criminal law and procedure.

Be it enacted by the General Assembly of the State of Indiana:

SECTION 1. IC 35-41-1-26.3, AS ADDED BY P.L.182-2011, SECTION 10, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2012]: Sec. 26.3. "Synthetic cannabinoid" means:

(1) a substance containing one (1) or more of the following chemical compounds, **including an analogue or homologue of the compound:**

~~(1)~~ **(A)** JWH-015 ((2-Methyl-1-propyl-1H-indol-3-yl)-1-naphthalenylmethanone).

~~(2)~~ **(B)** JWH-018 (1-pentyl-3-(1-naphthoyl)indole).

~~(3)~~ **(C)** JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).

~~(4)~~ **(D)** JWH-073 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).

~~(5)~~ **(E)** JWH-081 (4-methoxynaphthalen- 1-yl-(1-pentylindol- 3-yl)methanone).

~~(6)~~ **(F)** JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).

~~(7) JWH-200 (1-(2-morpholin-4-ylethyl)indol-3-yl)-~~

~~naphthalen-1-ylmethanone).~~
(G) JWH-200 ((1-(2-morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-ylmethanone).
~~(8)~~ **(H)** JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).
~~(9)~~ **(I)** JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).
~~(10)~~ **(J)** JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).
~~(11)~~ **(K)** HU-210 ((6aR,10aR)- 9-(Hydroxymethyl)-6,6-dimethyl- 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo [c]chromen- 1-ol).
~~(12)~~ **(L)** HU-211 ((6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl- 3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).
~~(13)~~ **(M)** HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-(2-methyloctan- 2-yl)phenyl]-7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).
~~(14) HU-331 ((3-hydroxy-2- [(1R,6R)-3-methyl-6-(1-methylethenyl)-2 -cyclohexen-1-yl]-5 -pentyl-2,5-cyclohexadiene-1,4-dione).~~
(N) HU-331 (3-hydroxy-2- [(1R,6R)-3-methyl-6-(1-methylethenyl)-2 -cyclohexen-1-yl]-5 -pentyl-2,5-cyclohexadiene-1,4-dione).
~~(15)~~ **(O)** CP 55,940 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]-5- (2-methyloctan-2-yl)phenol).
~~(16)~~ **(P)** CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]- 5-(2-methyloctan-2-yl)phenol) and its homologues.
~~(17)~~ **(Q)** WIN 55212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl) pyrrolo [1,2,3-de)- 1,4- benzoxazin-6-yl]-1-napthalenylmethanone).
~~(18)~~ **(R)** RCS-4 ((4-methoxyphenyl) (1-pentyl-1H-indol-3-yl)methanone).
~~(19)~~ **(S)** RCS-8 (1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl)ethanone).
~~(20)~~ **(T)** 4-Methylmethcathinone. Other name: mephedrone.
~~(21)~~ **(U)** 3,4-Methylenedioxymethcathinone. Other name: methylone.
~~(22)~~ **(V)** Fluoromethcathinone.
~~(23)~~ **(W)** 4-Methoxymethcathinone. Other name: methedrone.

~~(24)~~ **(X)** 4-Ethylmethcathinone (4-EMC).
~~(25)~~ **(Y)** Methylenedioxypyrovalerone. Other name: MDPV.
(Z) JWH-007, or
1-pentyl-2-methyl-3-(1-naphthoyl)indole.
(AA) JWH-098, or
1-pentyl-2-methyl-3-(4-methoxy-1-naphthoyl)indole.
(BB) JWH-164, or
1-pentyl-3-(7-methoxy-1-naphthoyl)indole.
(CC) JWH-210, or
1-pentyl-3-(4-ethyl-1-naphthoyl)indole.
(DD) JWH-201, or
1-pentyl-3-(4-methoxyphenylacetyl)indole.
(EE) JWH-203, or
1-pentyl-3-(2-chlorophenylacetyl)indole.
(FF) AM-694, or
1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole.
(GG) CP 50,556-1, or
[(6S,6aR,9R,10aR)-9-hydroxy-6-methyl-3-[(2R)-5-phenyl pentan-2-yl]oxy-5,6,6a,7,8,9,10,10a-octahydrophenanthri din-1-yl] acetate.
(HH) Dimethylheptylpyran, or DMHP.
(II) 4-Methyl-alpha-pyrrolidinobutiophenone, or MPBP.
(2) Any compound structurally derived from 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane by substitution at the nitrogen atom of the indole ring by alkyl, haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethy l,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 naphthyl ring to any extent.
(3) Any compound structurally derived from 3-(1-naphthoyl)173 pyrrole by substitution at the nitrogen atom of the pyrrole 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 pyrrole ring to any extent and whether or not substituted in the naphthyl ring to any extent.
(4) Any compound structurally derived from 1-(1-naphthylmethyl)indene by substitution at the 3-position of the indene 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 indene ring to any extent and whether or not substituted in the naphthyl ring to any extent.

(5) 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 and whether or not substituted in the phenyl ring to any extent.

(6) 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.

(7) 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.

SECTION 2. IC 35-48-2-4, AS AMENDED BY P.L.182-2011, SECTION 12, IS AMENDED TO READ AS FOLLOWS [EFFECTIVE JULY 1, 2012]: Sec. 4. (a) The controlled substances listed in this section are included in schedule I.

(b) Opiates. Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted by rule of the board or unless listed in another schedule, whenever the existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation:

Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide) (9815)

Acetylmethadol (9601)

Allylprodine (9602)

Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9832)

Alphacetylmethadol (9603)
Alphameprodine (9604)
Alphamethadol (9605)
Alphamethylfentanyl (9814)
Benzethidine (9606)
Beta-hydroxy-3-methylfentanyl (9831). Other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide
Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide) (9830)
Betacetylmethadol (9607)
Betameprodine (9608)
Betamethadol (9609)
Betaprodine (9611)
Clonitazene (9612)
Dextromoramide (9613)
Diampromide (9615)
Diethylthiambutene (9616)
Difenoxin (9168)
Dimenoxadol (9617)
Dimepheptanol (9618)
Dimethylthiambutene (9619)
Dioxaphetyl butyrate (9621)
Dipipanone (9622)
Ethylmethylthiambutene (9623)
Etonitazene (9624)
Etoxeridine (9625)
Furethidine (9626)
Hydroxypethidine (9627)
Ketobemidone (9628)
Levomoramide (9629)
Levophenacylmorphan (9631)
3-Methylfentanyl [N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenyl-propanimide](9813)
3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide) (9833)
MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) (9961)
Morpheridine (9632)
N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), including any isomers, salts, or salts of isomers (9818)
N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide

(thenylfentanyl), including any isomers, salts, or salts of isomers (9834)
Noracymethadol (9633)
Norlevorphanol (9634)
Normethadone (9635)
Norpipanone (9636)
Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide (9812)
Phenadoxone (9637)
Phenampromide (9638)
Phenomorphan (9647)
Phenoperidine (9641)
PEPAP [1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine] (9663)
Piritramide (9642)
Proheptazine (9643)
Properidine (9644)
Propiram (9649)
Racemoramide (9645)
Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide) (9835)
Tilidine (9750)
Trimeperidine (9646)

(c) Opium derivatives. Any of the following opium derivatives, their salts, isomers, and salts of isomers, unless specifically excepted by rule of the board or unless listed in another schedule, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

Acetorphine (9319)
Acetyldihydrocodeine (9051)
Benzylmorphine (9052)
Codeine methylbromide (9070)
Codeine-N-Oxide (9053)
Cyprenorphine (9054)
Desomorphine (9055)
Dihydromorphine (9145)
Drotebanol (9335)
Etorphine (except hydrochloride salt) (9056)
Heroin (9200)
Hydromorphinol (9301)
Methyldesorphine (9302)
Methyldihydromorphine (9304)
Morphine methylbromide (9305)

Morphine methylsulfonate (9306)
Morphine-N-Oxide (9307)
Myrophine (9308)
Nicocodeine (9309)
Nicomorphine (9312)
Normorphine (9313)
Pholcodine (9314)
Thebacon (9315)

(d) Hallucinogenic substances. Any material, compound, mixture, or preparation which contains any quantity of the following hallucinogenic, psychedelic, or psychogenic substances, their salts, isomers, and salts of isomers, unless specifically excepted by rule of the board or unless listed in another schedule, whenever the existence of these salts, isomers, and salts of isomers is possible within the specific chemical designation:

(1) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (7473). Other name: TCPy.
(2) 4-Bromo-2, 5-Dimethoxyamphetamine (7391). Some trade or other names: 4-Bromo-2, 5-Dimethoxy-a-methylphenethylamine; 4-Bromo-2, 5-DMA.
(3) 4-Bromo-2, 5-dimethoxphenethylamine (7392). Some trade or other names: 2-[4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B, Nexus.
(4) 2, 5-Dimethoxy-4-ethylamphet-amine (7399). Other name: DOET.
(5) 2, 5-Dimethoxy-4-(n)-propylthiophenethylamine (7348). Other name: 2C-T-7.
(6) 2, 5-Dimethoxyamphetamine (7396). Some trade or other names: 2, 5-Dimethoxy-a-methylphenethylamine; 2, 5-DMA.
(7) 4-Methoxyamphetamine (7411). Some trade or other names: 4-Methoxy-a-methylphenethylamine; Paramethoxyamphetamine; PMA.
(8) 5-Methoxy-3, 4-methylenedioxy amphetamine (7401). Other Name: MMDA.
(9) 5-Methoxy-N, N-diisopropyltryptamine, including any isomers, salts, or salts of isomers (7439). Other name: 5-MeO-DIPT.
(10) 4-methyl-2, 5-dimethoxyamphetamine (7395). Some trade and other names: 4-methyl-2, 5-dimethoxy-a-methylphenethylamine; DOM; and STP.
(11) 3, 4-methylenedioxy amphetamine (7400). Other name:

MDA.
(12) 3,4-methylenedioxy-N-ethylamphetamine (7404). Other names: N-ethyl-alpha-methyl-3,4(methylenedioxy) phenethylamine; N-ethyl MDA; MDE; and MDEA.
(13) 3, 4-methylenedioxymethamphetamine (MDMA) (7405).
(14) 3, 4, 5-trimethoxy amphetamine (7390). Other name: TMA.
(15) Alpha-ethyltryptamine (7249). Some trade and other names: Etryptamine; Monase; [alpha]-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; [alpha]-ET; and AET.
(16) Alpha-methyltryptamine (7432). Other name: AMT.
(17) Bufotenine (7433). Some trade and other names: 3-(B-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminonethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N, N-dimethyltryptamine; mappine.
(18) Diethyltryptamine (7434). Some trade or other names: N, N-Diethyltryptamine; DET.
(19) Dimethyltrytamine (7435). Some trade or other names: DMT.
(20) Ibogaine (7260). Some trade and other names: 7-Ethyl-6, 6b, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6, 9-methano-5H-pyrido (1', 2': 1, 2, azepino 4, 5-b) indole; tabernanthe iboga.
(21) Lysergic acid diethylamide (7315). Other name: LSD.
(22) Marijuana (7360).
(23) Mescaline (7381).
(24) Parahexyl (7374). Some trade or other names: 3-Hexyl-1-hydroxy-7, 8, 9, 10-Tetrahydro-6, 6, 9-trimethyl-6H-dibenzo (b,d) pyran; Snyhexyl.
(25) Peyote (7415), including:
    (A) all parts of the plant that are classified botanically as lophophora williamsii lemaire, whether growing or not;
    (B) the seeds thereof;
    (C) any extract from any part of the plant; and
    (D) every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or extracts.
(26) N-ethyl-3-piperidyl benzilate (7482). Other name: DMZ.
(27) N-hydroxy-3,4-methylenedioxyamphetamine (7402). Other names: N-hydroxy-alpha-methyl-3,4 (methylenedioxy)phenethylamine; and N-hydroxy MDA.
(28) N-methyl-3-piperidyl benzilate (7484). Other name: LBJ.

(29) Psilocybin (7437).
(30) Psilocyn (7438).
(31) Tetrahydrocannabinols (7370), including synthetic equivalents of the substances contained in the plant, or in the resinous extractives of Cannabis, sp. and synthetic substances, derivatives, and their isomers with similar chemical structure and pharmacological activity such as:
(A) π^{1} cis or trans tetrahydrocannabinol, and their optical isomers;
(B) π^{6} cis or trans tetrahydrocannabinol, and their optical isomers; and
(C) $\pi^{3,4}$ cis or trans tetrahydrocannabinol, and their optical isomers.
Since nomenclature of these substances is not internationally standardized, compounds of these structures, regardless of numerical designation of atomic positions are covered. Other name: THC.
(32) Ethylamine analog of phencyclidine (7455). Some trade or other names: N-Ethyl-1-phenylcyclohexylamine; (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl) ethylamine; cyclohexamine; PCE.
(33) Pyrrolidine analog of phencyclidine (7458). Some trade or other names: 1-(1-phenylcyclohexyl)-pyrrolidine; PCP_{y}; PHP.
(34) Thiophene analog of phencyclidine (7470). Some trade or other names: 1-(1-(2-thienyl) cyclohexyl) piperidine; 2-Thienyl Analog of Phencyclidine; TPCP.
(35) Synthetic cannabinoids **(as defined in IC 35-41-1-26.3).** ~~including a substance containing one (1) or more of the following chemical compounds:~~
~~(A) JWH-015 ((2-Methyl-1-propyl-1H-indol-3-yl)-1-naphthalenylmethanone).~~
~~(B) JWH-018 (1-pentyl-3-(1-naphthoyl)indole).~~
~~(C) JWH-019 (1-hexyl-3-(naphthalen-1-oyl)indole).~~
~~(D) JWH-073 (naphthalen-1-yl-(1-butylindol-3-yl)methanone).~~
~~(E) JWH-081 (4-methoxynaphthalen- 1-yl- (1-pentylindol-3-yl)methanone).~~
~~(F) JWH-122 (1-Pentyl-3-(4-methyl-1-naphthoyl)indole).~~
~~(G) JWH-200 (1-(2-morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-ylmethanone).~~
~~(H) JWH-250 (1-pentyl-3-(2-methoxyphenylacetyl)indole).~~
~~(I) JWH-251 (1-pentyl-3-(2-methylphenylacetyl)indole).~~

~~(J) JWH-398 (1-pentyl-3-(4-chloro-1-naphthoyl)indole).~~
~~(K) HU-210 ((6aR,10aR)- 9-(Hydroxymethyl)- 6,6-dimethyl- 3-(2-methyloctan-2-yl)- 6a,7,10,10a-tetrahydrobenzo [c]chromen- 1-ol).~~
~~(L) HU-211 ((6aS,10aS)-9-(Hydroxymethyl)- 6,6-dimethyl-3-(2-methyloctan-2-yl)- 6a,7,10,10a-tetrahydrobenzo [c]chromen-1-ol).~~
~~(M) HU-308 ([(1R,2R,5R)-2-[2,6-dimethoxy-4-(2-methyloctan- 2-yl)phenyl]- 7,7-dimethyl-4-bicyclo[3.1.1]hept-3-enyl] methanol).~~
~~(N) HU-331 ((3-hydroxy-2- [(1R,6R)-3-methyl-6-(1-methylethenyl)-2 -cyclohexen-1-yl]-5 -pentyl-2,5-cyclohexadiene-1,4-dione).~~
~~(O) CP 55,940 (2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl) cyclohexyl]- 5-(2-methyloctan-2-yl)phenol).~~
~~(P) CP 47,497 (2-[(1R,3S)-3-hydroxycyclohexyl]- 5-(2-methyloctan-2-yl)phenol) and its homologues.~~
~~(Q) WIN 55212-2 ((R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl) pyrrolo [1,2,3-de)- 1,4- benzoxazin-6-yl]-1-napthalenylmethanone).~~
~~(R) RCS-4 ((4-methoxyphenyl) (1-pentyl-1H-indol-3-yl)methanone).~~
~~(S) RCS-8 (1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl)ethanone).~~
~~(T) 4-Methylmethcathinone. Other name: mephedrone.~~
~~(U) 3,4-Methylenedioxymethcathinone. Other name: methylone.~~
~~(V) Fluoromethcathinone.~~
~~(W) 4-Methoxymethcathinone. Other name: methedrone.~~
~~(X) 4-Ethylmethcathinone. Other name: 4-EMC.~~
~~(Y) Methylenedioxypyrovalerone. Other name: MDPV.~~

(36) Salvia divinorum or salvinorin A, including:

(A) all parts of the plant that are classified botanically as salvia divinorum, whether growing or not;

(B) the seeds of the plant;

(C) any extract from any part of the plant; and

(D) every compound, manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or extracts.

(e) Depressants. Unless specifically excepted in a rule adopted by the board 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:

Gamma-hydroxybutyric acid (other names include GHB; gamma-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate) (2010)
Mecloqualone (2572)
Methaqualone (2565)

(f) Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation that contains any quantity of the following substances having a stimulant effect on the central nervous system, including its salts, isomers, and salts of isomers:

([+/-]) cis-4-methylaminorex (([+/-])cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine) (1590)
Aminorex (1585). Other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-dihydro-5-phenyl-2-oxazolamine.
Cathinone (1235). Some trade or other names: 2-amino-1-phenyl-1-propanone; alpha-aminopropiophenone; 2-aminopropiophenone; and norephedrone.
Fenethylline (1503).
N-Benzylpiperazine (7493). Other names: BZP; and 1-benzylpiperazine.
N-ethylamphetamine (1475)
Methcathinone (1237) Some other trade names: 2-Methylamino-1-Phenylpropan-I-one; Ephedrone; Monomethylpropion; UR 1431.
N, N-dimethylamphetamine (1480). Other names: N, N-alpha-trimethyl-benzeneethanamine; and N, N-alpha-trimethylphenethylamine.