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3 STUTTS AMENDMENT TO HB2, AS ENGROSSED
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8 On page 1, line 10, after "substances;" insert the
9 following:

10 to amend Section 20-2-23, Code of Alabama 1975,
11 relating to Schedule I of the controlled substances list, to
12 include additional controlled substances;
13

14 On page 2, delete lines 16 through 20.
15

16 On page 3, after line 21, insert the following
17 section and renumber the remaining sections accordingly:

18 "Section 2. Section 20-2-23, Code of Alabama 1975,
19 is amended to read as follows:

20 §Section 20-2-23.

21 (a) The Legislature finds the following:

22 (1) New synthetic substances are being created which
23 are not controlled under the provisions of existing state law
24 but which have a potential for abuse similar to or greater
25 than that for substances controlled under existing state law.
26 These new synthetic substances are called "synthetic
27 controlled substances or synthetic controlled substance

1 analogues" and can be designed to produce a desired
2 pharmacological effect and to evade the controlling statutory
3 provisions. Synthetic controlled substances or synthetic
4 controlled substance analogues are being manufactured,
5 distributed, possessed, and used as substitutes for controlled
6 substances.

7 (2) The hazards attributable to the traffic in and
8 use of a synthetic controlled substance or synthetic
9 controlled substance analogues are increased because their
10 unregulated manufacture produces variations in purity and
11 concentration.

12 (3) Many new synthetic substances are untested, and
13 it cannot be immediately determined whether they have useful
14 medical or chemical purposes.

15 (4) The uncontrolled importation, manufacture,
16 distribution, possession, or use of controlled substance
17 analogues has a substantial and detrimental impact on the
18 health and safety of the people of this state.

19 (5) Synthetic controlled substances or synthetic
20 controlled substance analogues can be created more rapidly
21 than they can be identified and controlled by action of the
22 Legislature. There is a need for a speedy determination of
23 their proper classification under existing law. It is
24 therefore necessary to identify and classify new substances
25 that have a potential for abuse, so that they can be
26 controlled in the same manner as other substances controlled
27 under existing state law.

1 (b) The controlled substances listed in this section
2 are included in Schedule I:

3 (1) Any of the following opiates, including their
4 isomers, esters, ethers, salts, and salts of isomers, esters
5 and ethers, unless specifically excepted, whenever the
6 existence of these isomers, esters, ethers and salts is
7 possible within the specific chemical designation:

- 8 a. Acetylmethadol;
- 9 b. Allylprodine;
- 10 c. Alphacetylmethadol;
- 11 d. Alphameprodine;
- 12 e. Alphamethadol;
- 13 f. Benzethidine;
- 14 g. Betacetylmethadol;
- 15 h. Betameprodine;
- 16 i. Betamethadol;
- 17 j. Betaprodine;
- 18 k. Clonitazene;
- 19 l. Dextromoramide;
- 20 m. Dextrophan;
- 21 n. Diampromide;
- 22 o. Diethylthiambutene;
- 23 p. Dimenoxadol;
- 24 q. Dimepheptanol;
- 25 r. Dimethylthiambutene;
- 26 s. Dioxaphetyl butyrate;
- 27 t. Dipipanone;

- 1 u. Ethylmethylthiambutene;
- 2 v. Etonitazene;
- 3 w. Etoxeridine;
- 4 x. Furethidine;
- 5 y. Hydroxypethidine;
- 6 z. Ketobemidone;
- 7 aa. Levomoramide;
- 8 bb. Levophenacylmorphan;
- 9 cc. Morpheridine;
- 10 dd. Noracymethadol;
- 11 ee. Norlevorphanol;
- 12 ff. Normethadone;
- 13 gg. Norpipanone;
- 14 hh. Phenadoxone;
- 15 ii. Phenampromide;
- 16 jj. Phenomorphan;
- 17 kk. Phenoperidine;
- 18 ll. Piritramide;
- 19 mm. Proheptazine;
- 20 nn. Properidine;
- 21 oo. Racemoramide;
- 22 pp. Trimeperidine.

23 (2) Any of the following opium derivatives, their
24 salts, isomers and salts of isomers, unless specifically
25 excepted, whenever the existence of these salts, isomers and
26 salts of isomers is possible within the specific chemical
27 designation:

- 1 a. Acetorphine;
- 2 b. Acetyldihydrocodeine;
- 3 c. Benzylmorphine;
- 4 d. Codeine methylbromide;
- 5 e. Codeine-N-Oxide;
- 6 f. Cyprenorphine;
- 7 g. Desomorphine;
- 8 h. Dihydromorphine;
- 9 i. Etorphine;
- 10 j. Heroin;
- 11 k. Hydromorphinol;
- 12 l. Methyldesorphine;
- 13 m. Methyldihydromorphine;
- 14 n. Morphine methylbromide;
- 15 o. Morphine methylsulfonate;
- 16 p. Morphine-N-Oxide;
- 17 q. Myrophine;
- 18 r. Nicocodeine;
- 19 s. Nicomorphine;
- 20 t. Normorphine;
- 21 u. Pholcodine;
- 22 v. Thebacon.

23 (3) Any material, compound, mixture or preparation
24 which contains any quantity of the following hallucinogenic
25 substances, their salts, isomers and salts of isomers, unless
26 specifically excepted, whenever the existence of these salts,

1 isomers and salts of isomers is possible within the specific
2 chemical designation:

- 3 a. 3,4-methylenedioxy amphetamine;
- 4 b. 5-methoxy-3,4-methylenedioxy amphetamine;
- 5 c. 3,4,5-trimethoxy amphetamine;
- 6 d. Bufotenine;
- 7 e. Diethyltryptamine;
- 8 f. Dimethyltryptamine;
- 9 g. 4-methyl-2,5-dimethoxy amphetamine;
- 10 h. Ibogaine;
- 11 i. Lysergic acid diethylamide;
- 12 j. Marihuana;
- 13 k. Mescaline;
- 14 l. Peyote;
- 15 m. N-ethyl-3-piperidyl benzilate;
- 16 n. N-methyl-3-piperidyl benzilate;
- 17 o. Psilocybin;
- 18 p. Psilocyn;
- 19 q. Tetrahydrocannabinols, except for

20 tetrahydrocannabinols in hemp, as defined in Section 2-8-381.

21 (4)a. A synthetic controlled substance that is any
22 material, mixture, or preparation that contains any quantity
23 of the following chemical compounds, their salts, isomers and
24 salts of isomers, unless specifically excepted, whenever the
25 existence of these salts, isomers and salts of isomers is
26 possible within the specific chemical designation or compound:

- 1 1. 3,4-Methylenedioxy-methcathinone (Methylone), some
2 trade or other names: 3,4-methylenedioxy-N-methylcathinone.
- 3 2. 3,4-Methylenedioxy-pyrovalerone, some other trade
4 names: (MDPV).
- 5 3. 4-Methylmethcathinone (Mephedrone), some trade or
6 other names: 4-methylephedrone.
- 7 4. 4-Methoxymethcathinone (Methedrone), some trade
8 or other names: bk-PMMA.
- 9 5. 3-Fluoromethcathinone, some trade or other names:
10 3-FMC.
- 11 6. 4-Fluoromethcathinone (Flephedrone), some trade
12 or other names: 4-FMC.
- 13 7.
14 1-[(5-fluoropentyl)-1H-indol-3-yl]-(2-iodophenyl)methanone,
15 some trade or other names: AM-694.
- 16 8.
17 1-[(5-fluoropentyl)-1H-indol-3-yl]-(naphthalen-1-yl)methanone,
18 some trade or other names: AM-2201.
- 19 9. (6aR, 10aR)-9-(hydroxymethyl)-6,
20 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[
21 c]chromen-1-ol, some trade or other names: HU-210.
- 22 10.
23 (6aS,10aS)-9-(Hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-y
24 l)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol, some trade or
25 other names: HU-211, Dexanabinol.
- 26 11. 1-Pentyl-2-methyl-3-(1-naphthoyl)indole, some
27 trade or other names: JWH-007.

1 12.
2 (2-Methyl-1-propyl-1H-indol-3-yl)-1-naphthalenylmethanone,
3 some trade or other names: JWH-015.
4 13. Naphthalen-1-yl-(1-pentylindol-3-yl)methanone,
5 some trade or other names: JWH-018.
6 14. 1-Hexyl-3-(naphthalen-1-oyl)indole, some trade
7 or other names: JWH-019.
8 15. Naphthalen-1-yl-(butylindol-3-yl)methanone, some
9 trade or other names: JWH-073.
10 16.
11 4-Methoxynaphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
12 trade or other names: JWH-081.
13 17.
14 4-Methoxynaphthalen-1-yl-(1-pentyl-2-methylindol-3-yl)
15 methanone, some trade or other names: JWH-098.
16 18.
17 4-Methylnaphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
18 trade or other names: JWH-122.
19 19.
20 (1-(2-Morpholin-4-ylethyl)indol-3-yl)-naphthalen-1-ylmethanone
21 , some trade or other names: JWH-200.
22 20.
23 2-(2-Chlorophenyl)-1-(1-pentylindol-3-yl)ethanone, some trade
24 or other names: JWH-203.
25 21.
26 4-Ethyl-naphthalen-1-yl-(1-pentylindol-3-yl)methanone, some
27 trade or other names: JWH-210.

1 22.
2 2-(2-Methoxyphenyl)-1-(1-pentylindol-3-yl)ethanone, some trade
3 or other names: JWH-250.
4 23.
5 5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-ylmethano
6 ne, some trade or other names: JWH-307.
7 24. 1-Pentyl-3-(4-Chloro-1-naphthoyl)indole, some
8 trade or other names: JWH-398.
9 25.
10 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol
11 (Cannabicyclohexanol), some trade or other names: CP 47, 497,
12 and homologues.
13 26.
14 2-(2-Methoxyphenyl)-1-[1-(2-cyclohexylethyl)indol-3-yl]ethanon
15 e, some trade or other names: RCS-8, SR-18.
16 27.
17 2-(4-Methoxyphenyl)-1-(1-pentyl-indol-3-yl)methanone, some
18 trade or other names: RCS-4.
19 28.
20 (R)-(+)-[2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1
21 ,2,3-de]-1,4-benzoxazin-6-yl]-1-naphthalenylmethanone, some
22 trade or other names: WIN 55,212-2.
23 29.
24 (4-Methoxyphenyl)-[2-methyl-1-(2-morpholin-4-ylethyl)indol-3-y
25 l]methanone, some trade or other names: WIN 48,098,
26 Pravadoline.

1 b. In addition to any material, mixture, or
2 preparation that contains any quantity of the chemical
3 compounds listed in paragraph a., a synthetic controlled
4 substance also includes the following chemical compounds,
5 their salts, isomers and salts of isomers, unless specifically
6 excepted, whenever the existence of these salts, isomers and
7 salts of isomers is possible within the specific chemical
8 designation or compound:

9 1.

10 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole,
11 some trade or other names: (AM-2233).

12 2. 1-Pentyl-3-(1-adamantoyl)indole, some trade or

13 other names: (AB001).

14 3.

15 [1-[(1-methyl-2-piperidinyl)methyl]-1H-indol-3-yl]-1-naphthale
16 nyl-methanone, some trade or other names: (AM1220).

17 4.

18 1-(5-Fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole,
19 some trade or other names: (XLR11).

20 5.

21 1-Pentyl-3-(2,2,3,3-tetramethylcyclopropoyl)indole, some trade
22 or other names: (UR-144).

23 6.

24 6-Methyl-2[(4-methylphenyl)amino]-4H-3,1-benzoxazin-4-one,
25 some trade or other names: (URB 754).

26 7. [1,1'-biphenyl]-3-yl-carbamic acid, cyclohexyl

27 ester, some trade or other names: (URB 602).

1 8.
2 (3'-(Aminocarbonyl)[1,1'-biphenyl]-3-yl)-cyclohexylcarbamate,
3 some trade or other names: (URB597).
4 9.
5 1-(5-Fluoropentyl)-3-(4-methyl-1-naphthoyl)indole, some trade
6 or other names: (MAM2201).
7 10.
8 1-naphthalenyl[4-(pentyloxy)-naphthalenyl]methanone, some
9 trade or other names: (CB-13).
10 11.
11 1-(5-Chloropentyl)-3-(2,2,3,3-tetramethylcyclopropoyl)indole,
12 some trade or other names: (5-Chloro-UR-144).
13 12.
14 1-(5-Fluoropentyl)-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indole-3-
15 carboxamide, some trade or other names: (STS-135).
16 13.
17 1[(N-Methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl)indole,
18 some trade or other names: (AM1248).
19 14. N-Adamantyl-1-pentyl-1H-indole-3-carboxamide,
20 some trade or other names: (SDB-001, 2NE1).
21 15.
22 1-Pentyl-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indazole-3-carboxam
23 ide, some trade or other names: (AKB48, APINACA).
24 16. 3-Naphthoylindole.
25 17.
26 1-[2-(4-Morpholinyl)ethyl]-3-(2,2,3,3-tetramethylcyclopropyl) i
27 ndole, some trade or other names: (A 796,260).

1 18.
2 1-[(tetrahydropyran-4-ylmethyl)-1H-indol-3-yl]-(2,2,3,3-tetram
3 ethylcyclopropyl)methanone, some trade or other names: (A
4 834,735).
5 19.
6 1-(Pent-4-en-1-yl)-3-(4-methyl-1-naphthoyl)indole, some trade
7 or other names: (JWH-122 4-pentenyl analog).
8 20.
9 N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-fluorophenyl)m
10 ethyl]-1H-indazole-3-carboxamide some trade or other names:
11 (AB-FUBINACA).
12 21.
13 [1-(5-bromopentyl)-1H-indol-3-yl]-(2,2,3,3-tetramethylcycloprop
14 yl)methanone, some trade or other names: (5-Bromo-UR-144)
15 22.
16 5-(1,1-Dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexylphenol,
17 some trade or other names: (CP-47,497 C8 homolog).
18 23.
19 1-(5-Fluoropentyl)-N-tricyclo[3,3,1,13,7]dec-1-yl-1H-indazole-
20 3-carboxamide, some trade or other names: (5F-AKB48,
21 5F-APINACA).
22 24. 1-(penta-4-ene)-3-(1-naphthoyl)indole, some
23 trade or other names: (JWH-022).
24 25. 1-(5-Chloropentyl)-3-(1-naphthoyl)indole, some
25 trade or other names: (Chloro-AM-2201, JWH-018
26 N-5-chloropentyl analog).

1 26. 1-(5-Hydroxypentyl)-3-(1-naphthoyl)indole, some
2 trade or other names: (Hydroxy-AM-2201).

3 27.
4 N-[(2E)-3-(2-Methoxyethyl)4,5-dimethyl-1,3-thiazole-2(3H)-ylid
5 ene]-2,2,3,3-tetramethylcyclopropane carboxamide, some trade
6 or other names: (A 836,339).

7 28. 1-Pentyl-3-(2-iodobenzoyl)indole, some trade or
8 other names: (AM 679).

9 29. 1-Pentyl-3-(2-methylphenacetyl)indole, some
10 trade or other names: (JWH-251).

11 30. 1-pentyl-1H-indole-3-carboxylic acid
12 8-quinolinyl ester, some trade or other names: (PB-22, QUPIC).

13 31. 1-(5-fluoropentyl)-1H-indole-3-carboxylic acid
14 8-quinolinyl ester, some trade or other names: (5F-PB-22).

15 32.
16 1-pentyl-N-(naphthalen-1-yl)-1H-indole-3-carboxamide, some
17 trade or other names: (MN-24, NNE1).

18 33. 1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid
19 8-quinolinyl ester, some trade or other names: (BB-22,
20 QUCHIC).

21 34.
22 N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-pentyl-1H-indazole
23 -3-carboxamide, some trade or other names: (AB-PINACA).

24 35.
25 7-methoxy-1-(2-morpholinoethyl)-N-((1S,2S,4R)-1,3,3-trimethylb

1 bicyclo[2.2.1]heptan-2-yl)-1H-indole-3-carboxamide, some trade
2 or other names: (MN-25).

3 36. ADB-PINACA.

4 37. FUB-AKB-48.

5 38. FUB-PB-22.

6 39. Heptyl-UR144.

7 40. THJ-018.

8 41. THJ-2201.

9 42. 1-heptyl-3-(1-naphthoyl)indole), some trade or
10 other names: (JWH-20).

11 43.

12 Napthalen-1-yl-(1-propyl-1H-indol-3-yl)methanone, some trade
13 or other names: (JWH-072).

14 44. (6aR,10aR)-3-(1, 1-Dimethylbutyl)-6a, 7, 10,
15 10a-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran, some
16 trade or other names: (JWH-133).

17 45. 3-(napthalen-1-ylmethyl)-1-pentyl-1H-indole,
18 some trade or other names: (JWH-175).

19 46. 1-pentyl-3-(4-methoxyphenylacetyl)indole, some
20 trade or other names: (JWH-201).

21 47. 1-pentyl-3-(3-methoxyphenylacetyl)indole, some
22 trade or other names: (JWH 302).

23 48.

24 [(1R,2R,5R)-2-[2,6-dimethoxy-4-(2-methyloctan-2-yl)phenyl]-7,7
25 -dimethyl-4-bicyclo[3.1.1]hept-3-enyl]methanol, some trade or
26 other names: (HU-308).

1 49.
2 3-hydroxy-2-[(1R,6R)-3-methyl-6-(1-methylethenyl)-2-cyclohexen
3 -1-yl]-5-pentyl-2,5-cyclohexadiene-1,4-dione, some trade or
4 other names: (HU-331).
5 50.
6 N-cyclopropyl-11-(3-hydroxy-5-pentylphenoxy)-undecanamide,
7 some trade or other names: (CB-25).
8 51.
9 N-cyclopropyl-11-(2-hexyl-5-hydroxyphenoxy)-undecanamide, some
10 trade or other names: (CB-52).
11 52.
12 2-[(1R,2R,5R)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-5-(2-me
13 thyloctan-2-yl)phenol, some trade or other names:
14 (CB-55,940) (CB-55).
15 53. 4-Methylethylcathinone, some trade or other
16 names: (4-MEC, 4-Methylethcathinone).
17 54. 4'-Methyl-alpha-pyrrolidinopropiophenone, some
18 trade or other names: (MPPP, ZZ-1).
19 55.
20 (RS)-1-naphthalen-2-yl-2-pyrrolidin-1-ylpentan-1-one, some
21 trade or other names: (Naphyrone).
22 56. alpha,alpha-Diphenyl-2-piperidinemethanol, some
23 trade or other names: (Pipradrol, Meratran).
24 57.
25 (RS)-1-(4-methylphenyl)-2-(1-pyrrolidinyl)pentan-1-one, some
26 trade or other names: (Pyrovalerone).

1 58. 3,4-Dimethylmethcathinone, some trade or other
2 names: (3,4-DMMC).

3 59. 4-Fluoroamphetamine, some trade or other names:
4 (4-FA).

5 60. 4-Fluoromethamphetamine, some trade or other
6 names: (4-FMA).

7 61. Butylone, some trade or other names: (bk-MBDB).

8 62. alpha-Pyrrolidinopentiophenone, some trade or
9 other names: (alpha-PVP).

10 63. beta-keto-Dimethylbenzodioxolylbutanamine, some
11 trade or other names: (bk-DMBDB).

12 64. 2-(methylamino)-1-phenylbutan-1-one, some trade
13 or other names: (Buphedrone).

14 65. (RS)-2-ethylamino-1-phenyl-propan-1-one, some
15 trade or other names: (N-Ethylcathinone).

16 66. 2-Fluoroamphetamine, some trade or other names:
17 (2-FA).

18 67. Methoxetamine, some trade or other names: (MXE).

19 68. 2-Methylamino-1-phenylpentan-1-one, some trade
20 or other names: (Pentedrone).

21 69. 3,4-Methylenedioxycathinone, some trade or other
22 names: (MDC).

23 70. 2-Fluoromethamphetamine, some trade or other
24 names: (2-FMA).

25 71. 4-methylmethamphetamine, some trade or other
26 names: (4-MMA).

1 72. 4-Fluoroisocathinone, some trade or other names:
2 (4-FIC).

3 73. 3-Fluoromethamphetamine, some trade or other
4 names: (3-FMA).

5 74. Methiopropamine, some trade or other names:
6 (MPA).

7 75. alpha-Pyrrolidinobutiophenone, some trade or
8 other names: (alpha-PBP).

9 76. 4-Methoxy-N-methylcathinone, some trade or other
10 names: (Methedrone, bk-PMMA).

11 77. alpha-Pyrrolidinopropiophenone, some trade or
12 other names: (alpha-PPP).

13 78. (RS)-2-benzhydrylpiperidine, some trade or other
14 names: (Desoxypipradrol).

15 79. 3,4-Methylenedioxyethylcathinone, some trade or
16 other names: (MDEC).

17 80.
18 3,4-Methylenedioxy-alpha-pyrrolidinobutiophenone, some trade
19 or other names: (MDPBP).

20 81.
21 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one
22 (Pentylone, bk-MBDP).

23 82. 3-Fluoroamphetamine, some trade or other names:
24 (3-FA).

25 83. 3-Fluoromethcathinone, some trade or other
26 names: (3-FMC).

1 84. 2-Fluoromethcathinone, some trade or other
2 names: (2-FMC).
3 85.
4 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)propan-1-one
5 (bk-MDDMA).
6 86. N,N-Diethylcathinone, some trade or other names:
7 (Amfepramone, DEC).
8 87. 1,3-Dimethylamylamine, some trade or other
9 names: (DMAA).
10 88. N, N-Dimethylcathinone, some trade or other
11 names: (DMC).
12 89. N-Ethyl-3,4-methylenedioxycathinone, some trade
13 or other names: (bk-MDEA).
14 90. N-Ethylamphetamine, some trade or other names:
15 (EMA).
16 91. N-Ethylcathinone, some trade or other names:
17 (EC).
18 92. 2-Ethylethcathinone, some trade or other names:
19 (2-EEC).
20 93. 4-Ethyl-N-ethylcathinone, some trade or other
21 names: (4-EEC).
22 94.
23 2-(5-Methoxy-1-benzofuran-3-yl)-N,N-dimethylethanamine, some
24 trade or other names: (Dimembfe).
25 95. 2-(5-Methoxy-1-benzofuran-3-yl)N-ethylethamine.
26 96. 4-Methoxymethamphetamine, some trade or other
27 names: (PMMA).

1 97. 4-Methoxy-N-ethylamphetamine, some trade or
2 other names: (PMEA).
3 98. 4-Methoxy-N-ethylcathinone, some trade or other
4 names: (ETHEDRONE).
5 99. 3-Methylmethcathinone, some trade or other
6 names: (3-MMC).
7 100. 4-Methyl-alpha-pyrrolidinobutiophenone, some
8 trade or other names: (MPBP).
9 101. 2-Methylethcathinone, some trade or other
10 names: (2-MEC).
11 102. 3-Methylethcathinone, some trade or other
12 names: (3-MEC).
13 103. 2-Ethylethcathinone, some trade or other names:
14 (2-EEC).
15 104. 3-Ethylethcathinone, some trade or other names:
16 (3-EEC).
17 105. 3-Ethylmethcathinone, some trade or other
18 names: (3-EMC).
19 106.
20 3',4'-Methylenedioxy-alpha-pyrrolidinopropiophenone, some
21 trade or other names: (MDPPP).
22 107. alpha-Pyrrolidinopentiothiophenone, some trade
23 or other names: (alpha-PVT).
24 108. 3-Methoxymethcathinone, some trade or other
25 names: (3-MeOMC).
26 109. N-Methyl-1,3-benzodioxolylbutanamine, some
27 trade or other names: (MBDB).

1 110. Ethcathinone, some trade or other names:
2 (ETHYLPROPION, ETH-CAT) .
3 111. Ethylone (3,4-methylenedioxy-N-ethylcathinone).
4 112. N-N-Diethyl-3,4-methylenedioxcathinone.
5 113. 3,4-methylenedioxy-propiofenone.
6 114. 2-Bromo-3,4-methylenedioxypropiofenone.
7 115. 3,4-methylenedioxy-propiofenone-2-oxime.
8 116. N-Acetyl-3,4-methylenedioxcathinone.
9 117. N-Acetyl-N-Methyl-3,4-methylenedioxcathinone.
10 118. N-Acetyl-N-Ethyl-3,4-methylenedioxcathinone.
11 119. 4-Bromomethcathinone.
12 120. 3-Bromomethcathinone.
13 121. Eutylone
14 (beta-Keto-Ethylbenzodioxolylbutanamine) .
15 122. 4'-Methoxy-alpha-pyrrolidinopropiofenone, some
16 trade or other names: (MOPPP) .
17 123. 4'-Methyl-alpha-pyrrolidinohexiofenone, some
18 trade or other names: (MPHP) .
19 124. Benocyclidine (BCP) or
20 Benzothiophenylcyclohexylpiperidine, some trade or other
21 names: (BTCP) .
22 125. 4-Fluoro-(methylamino)butyrofenone, some trade
23 or other names: (F-MABP) .
24 126. 3-Methyl-4-Methoxymethacathinone, some trade or
25 other names: (3-Me-4-MeO-MCAT) .

1 127. 4-Methyl-(ethylamino)-butryophenone, some trade
2 or other names: (Me-EABP).
3 128. 4-Ethyl-methcathinone, some trade or other
4 names: (4-EMC).
5 129. 4-methoxy-N-ethylcathinone
6 (bk-PMC;p-methox-ethcathinone).
7 130. 4'-Methoxy-alpha-pyrroldino-propiofenone
8 (MeOPPP; 4'-MeO-PPP).
9 131. 3-Fluorocathinone (3-FC).
10 132. 4-Fluorocathinone (4-FC).
11 133. 4-methyl-buphedrone (4-MeMABP; 4MeBP; BZ-6378).

12 134. 3,4-Methylenedioxy-N-benzylcathinone, some
13 trade or other names: (BMDP).
14 135. N-Benzyl-butylone, some trade or other names:
15 (BMDB).
16 136. N-Hydroxy-3,4-methlyenedioxymethcathinone.
17 137. N-ethylbuphedrone, some trade or other names:
18 (NEB).
19 138. 4-Fluorobuphedrone, some trade or other names:
20 (4-FBP).
21 139. 4-Methoxy-pyrrolidinbutrophenone (4-MeO-PBP).
22 140. 4-Ethyl-pyrrolidinobutrophenone, some trade or
23 other names: (4-Et-PBP).
24 141. 5-(2-aminopropyl)indole, some trade or other
25 names: (5-IT).
26 142. 1-phenyl-2-(piperidin-1-yl)butan-1-one.

- 1 143. 2,4,5-Trimethyl-methacathinone, some trade or
2 other names: (2,4,5-TMMC).
- 3 144. alpha-pyrrolidino-heptiophenone, some trade or
4 other names: (alpha-PHpP).
- 5 145. 4-Methylamphetamine (4-MA: pTAP; PAL-313;
6 4-MeA; PmeA).
- 7 146. N-Ethyl-methamphetamine.
- 8 147. 4-(2-Aminopropyl)benzofuran, some trade or
9 other names: (4-APB).
- 10 148. 5-(2-Aminopropyl)-2,3-dihydro-1H-indene
11 (5-APDI; IAP; AIP; indanylaminoporpane).
- 12 149. 6,7-Methylenedioxy-2-aminotetralin, some trade
13 or other names: (MDAT).
- 14 150. 4-Methylthioamphetamine (4-MTA; P1882).
- 15 151. 4-Chloroamphetamine (p-chloro-amphetamine).
- 16 152. 2,4,6-Trimethoxyamphetamine, some trade or
17 other names: (TMA-6).
- 18 153. 2,4,5-Trimethoxyamphetamine, some trade or
19 other names: (TMA-2).
- 20 154. 2,5-Dimethylamphetamine, some trade or other
21 names: (2,5-DMA).
- 22 155. 3,4-Dimethylamphetamine, some trade or other
23 names: (3,4-DMA).
- 24 156. N-propylamphetamine.
- 25 157. 4-Hydroxyamphetamine.
- 26 158. 3-Hydroxyamphetamine.

1 159. Methylenedioxydimethylamphetamine, some trade
2 or other names: (MDDM).
3 160. 2-Aminoindane, some trade or other names:
4 (2-AI).
5 161. 5,6-Methylenedioxy-N-methyl-aminoindane, some
6 trade or other names: (MDMAI).
7 162. 2C-T-21.
8 163. 2C-B-Fly.
9 164. 3,4-dimethyl-2,5-dimethoxyphenethylamine
10 (2C-G).
11 165. 25D-NBOMe.
12 166. 25G-NBOMe.
13 167. 25N-NBOMe.
14 168. Bromo-benzylidifuranyl-isopropylamine, some
15 trade or other names: (Bromo Dragon Fly).
16 169. 3C-B fly.
17 170. 2,5-Dimethoxy-4-ethylthioamphetamine, some
18 trade or other names: (Aleph-2).
19 171.
20 1-[(4-ethoxy-2,5-dimethoxy)phenyl]propan-2-amine, some trade
21 or other names: (MEM).
22 172.
23 1-[2,5-dimethoxy-4-(propylthio)phenyl]propan-2-amine, some
24 trade or other names: (Aleph-7).
25 173. N-benzyl-2-phenylethylamine.
26 174. N,N-dimethyl-2-phenylethylamine.

1 175. 6-chloro-2-aminotetralin, some trade or other
2 names: (6-CAT).
3 176. 2-phenylpropan-1-amine, some trade or other
4 names: (B-Me-PEA).
5 177. 2-Phenethylamine, some trade or other names:
6 (2-PEA).
7 178.
8 1-methylamino-1-(3,4-methylenedioxyphenyl)propane, some trade
9 or other names: (M-ALPHA).
10 179. Camfetamine.
11 180. Methoxyphenamine.
12 181. 4-methylaminorex, some trade or other names:
13 (4-MAR; 4-MAX; U4Euh; Euphoria; Ice).
14 182. (1-thiophen-2-yl)propan-2-amine
15 (Thienoamphetamine).
16 183. Dimethocaine.
17 184. 4-Fluoroephedrine.
18 185. 4-methyaminorex (p-methyl derivative).
19 186.
20 1-[(N-methylpiperidin-2-yl)methyl]-2-methyl-3-(naphthalen-1-oy
21 l)-6-nitroindole (AM1221).
22 187.
23 (1-butyl-1H-indol-3-yl)(4-methoxyphenyl)-methanone (RCS-4 (C4)
24 homolog).
25 188.
26 5-[3-(1-naphthoyl)-1H-indole-1-yl]pentanenitrile, some trade
27 or other names: (AM2232).

1 189. 1-(Pentyl)-3-(4-bromo-1-naphthoyl)-indole, some
2 trade or other names: (JWH-387).
3 190. 1-(Pentyl)-3-(4-fluoro-1-naphthoyl)-indole,
4 some trade or other names: (JWH-412).
5 191. 1-(5-chloropentyl)-3-(2-iodobenzoyl)indole, some
6 trade or other names: (AM694 Derivative).
7 192.
8 (2-iodo-5-nitrophenyl)-[1-[(1-methylpiperidin-2-yl)methyl]1H-i
9 ndol-3-yl]-methanone, some trade or other names: (AM1241).
10 193. 1-Pentyl-3-[1-(4-propyl)naphthoyl]indole, some
11 trade or other names: (JWH-182).
12 194. JWH-081 2-methoxynaphthyl isomer, some trade or
13 other names: (JWH-267).
14 195.
15 (3-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone, some trade
16 or other names: (RCS-4 3-methoxy isomer).
17 196.
18 [1-(5-fluoropentyl)-1H-indol-3-yl](4-ethyl-1-naphthalenyl)-met
19 hanone (EAM-2201).
20 197. ADB-FUBINACA.
21 198. ADBICA.
22 199. AM-279.
23 200. JWH-370.
24 201. NNE-1.
25 202. MAM-2201 chloropentyl derivative.
26 203. 1-(5-fluoropentyl)-3-(2-methyl-benzoyl)indole.
27 204. 1-(5-fluoropentyl)-3-(2-ethylbenzoyl)indole.

1 205. AB-005.
2 206. AB-005 Azepane isomer.
3 207.
4 4-hydroxy-3,3,4-trimethyl-1-(1-pentyl-1H-indol-3-yl)pentan-1-o
5 ne (4-HTMPIPO).
6 208. UR-12.
7 209. 5-Fluoro-ADBICA.
8 210. BAY-38-7271; KN 38-7271.
9 211. JTE-907.
10 212. Org 27569.
11 213. Org 27759.
12 214. Org 29647.
13 215. LY 2183240.
14 216. JTE 7-31.
15 217. URB 937.
16 218. 3-methoxy-eticyclidine, some trade or other
17 names: (3-MeO-PCE).
18 219. 1-Phenylcyclohexanamine, some trade or other
19 names: (PCA).
20 220. 4-Methyl-phencyclidine, some trade or other
21 names: (4-Me-PCP).
22 221. 4-Methoxy-eticyclidine, some trade or other
23 names: (4-MeO-PCE).
24 222. 4-Methoxyphencyclidine, some trade or other
25 names: (Methoxydine; 4MeO-PCP).
26 223. 3-Methoxyphencyclidine, some trade or other
27 names: (3-MeO-PCP).

1 224. 1-phenyl-N-propylcyclohexanamine, some trade or
2 other names: (PCPr).
3 225. N-(2-methoxyethyl)-1-phenylcyclohexanamine,
4 some trade or other names: (PCMEA).
5 226. N-(2-ethoxyethyl)-1-phenylcyclohexanamine, some
6 trade or other names: (PCEEA).
7 227. N-(3-methoxypropyl)-1-phenylcyclohexanamine,
8 some trade or other names: (PCMPA).
9 228. 3-Hydroxy-phencyclidine, some trade or other
10 names: (3-OH-PCP).
11 229. Methoxyketamine, some trade or other names:
12 (2-MeO-2-deschloro-ketamine).
13 230. Tiletamine, some trade or other names: (TCE).
14 231. N-ethylnorketamine.
15 232. N-Methyltryptamine, some trade or other names:
16 (NMT).
17 233. N-Methyl-N-isopropyltryptamine, some trade or
18 other names: (MiPT; MIPT).
19 234. 4-hydroxy-N,N-methylisopropyltryptamine, some
20 trade or other names: (4-OH-MiPT).
21 235. 4-Acetoxy-N,N-diisopropyl-tryptamine
22 (4-AcO-DiPT: 4-AcO-DIPT; 4-Acetoxy-MiPT).
23 236. 4-Methoxy-N,N-dimethyltryptamine, some trade or
24 other names: (4-MeO-DMT).
25 237. 5-Hydroxytryptamine, some trade or other names:
26 (5-HT).

1 238. 5-acetoxy-N,N-dimethyltryptamine, some trade or
2 other names: (5-AcO-DMT).
3 239. 5-Methoxy-N,N-dipropyltryptamine, some trade or
4 other names: (5-MeO-DPT).
5 240. d-Lysergic acid amide, some trade or other
6 names: (LSA; ergine).
7 241. 2,5-dimethoxy-4-chloroamphetamine, some trade
8 or other names: (DOC).
9 242.
10 N-(2-Methoxybenzyl)-4-iodo-2,5-dimethoxyphenethylamine, some
11 trade or other names: (25I-NBOMe).
12 243. 4-Ethyl-2,5-dimethoxyphenethylamine, some trade
13 or other names: (2C-E).
14 244. 2,5-Dimethoxy-4-iodophenethylamine, some trade
15 or other names: (2C-I).
16 245. 6-(2-Aminopropyl)-2,3-dihydrobenzofuran, some
17 trade or other names: (6-APDB).
18 246. 6-(2-Aminopropyl)benzofuran, some trade or
19 other names: (6-APB).
20 247. 5-(2-Aminopropyl)-2,3-dihydrobenzofuran, some
21 trade or other names: (5-APDB).
22 248. 5-(2-Aminopropyl)benzofuran, some trade or
23 other names: (5-APB).
24 249. 2,5-Dimethoxy-4-(n)-propylthiophenethylamine,
25 some trade or other names: (2C-T-7).
26 250. 2,5-Dimethoxy-4-(n)-propylphenethylamine, some
27 trade or other names: (2C-P).

1 251. 2,5-Dimethoxy-4-bromoamphetamine, some trade or
2 other names: (DOB).

3 252. 2,5-Dimethoxy-4-bromobenzylpiperazine, some
4 trade or other names: (2C-B-BZP).

5 253. 2,5-Dimethoxy-4-bromophenethylamine, some trade
6 or other names: (2C-B).

7 254. 2,5-Dimethoxy-4-chlorophenethylamine, some
8 trade or other names: (2C-C).

9 255. 2,5-Dimethoxy-(4-ethylthio)phenethylamine, some
10 trade or other names: (2C-T-2).

11 256. 2,5-Dimethoxy-4-iodoamphetamine, some trade or
12 other names: (DOI).

13 257. 2,5-Dimethoxy-4-methylamphetamine, some trade
14 or other names: (DOM).

15 258. 2,5-Dimethoxyphenethylamine, some trade or
16 other names: (2C-H).

17 259.

18 2-(2,5-Dimethoxyphenyl-4-bromo)-N-(2-methoxybenzyl)ethanamine,
19 some trade or other names: (25B-NBOMe).

20 260.

21 2-(2,5-Dimethoxyphenyl-4-chloro)-N-(2-methoxybenzyl)ethanamine
22 , some trade or other names: (25C-NBOMe).

23 261.

24 2-(2,5-Dimethoxyphenyl-4-ethyl)-N-(2-methoxybenzyl)ethanamine,
25 some trade or other names: (25E-NBOMe).

26 262. 2-Ethylmethcathinone, some trade or other
27 names: (2-EMC).

1 263.
2 2-(2,5-Dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine, some
3 trade or other names: (25H-NBOMe).
4 264. BZP (Benzylpiperazine).
5 265. para-Fluorophenylpiperazine.
6 266. 1-(4-Methylphenyl)piperazine.
7 267. meta-Cholorophenylpiperazine.
8 268. para-Methoxyphenylpiperazine.
9 269. DBZP (1,4-dibenzylpiperazine).
10 270. TFMPP (3-Trifluoromethylphenylpiperazine).
11 271. 2C-T-4
12 (2,5-Dimethoxy-4-isopropylthiophenethylamine).
13 272. 2C-T
14 (2,5-Dimethoxy-4-methylthiophenethylamine).
15 273. 2C-D
16 (2-(2,5-Dimethoxy-4-methylphenyl)ethanamine).
17 274. 2C-N 2,5-Dimethoxy-4-nitrophenethylamine.
18 275. 5-methoxy-N,N-diallyltryptamine, some trade or
19 other names: (5-MeO-DALT).
20 276. 5-Methoxy-N,N-Diisopropyltryptamine, some trade
21 or other names: (5-MeO-DIPT).
22 277. 5-Methoxy-alpha-methyltryptamine, some trade or
23 other names: (5-MeO-AMT).
24 278. 4-Acetoxy-N,N-dimethyltryptamine, some trade or
25 other names: (4-AcO-DMT).
26 279. 4-Hydroxy-N,N-diethyltryptamine, some trade or
27 other names: (4-HO-DET).

1 280. 4-Hydroxy-N,N-diisopropyltryptamine, some trade
2 or other names: (4-HO-DIPT).

3 281. 4-Hydroxy-N-methyl-N-ethyltryptamine, some
4 trade or other names: (4-OH-MET).

5 282. 5-Methoxy-N,N-diethyltryptamine, some trade or
6 other names: (5-MeO-DET).

7 283. 5-Methoxy-N-methyl-N-isopropyltryptamine, some
8 trade or other names: (5-MeO-MIPT).

9 284. 4-Acetoxy-N,N-diethyltryptamine, some trade or
10 other names: (4-AcO-DET).

11 285. 4-Acetoxy-N-methyl-N-isopropyltryptamine, some
12 trade or other names: (4-AcO-MIPT).

13 286. N,N-Dipropyltryptamine, some trade or other
14 names: (DPT).

15 287. N,N-Diisopropyltryptamine, some trade or other
16 names: (DIPT).

17 288. 4-Methoxy-N-methyl-N-isopropyltryptamine, some
18 trade or other names: (4-MeO-MIPT).

19 289. Tyramine (4-Hydroxyphenethylamine).

20 290. 5-Hydroxy-alpha-methyltryptamine.

21 291. 5-Hydroxy-N-methyltryptamine.

22 292. 5-Methoxy-N,N-dimethyltryptamine.

23 293. 5-Methyl-N,N-dimethyltryptamine.

24 294. Diphenylprolinol, some trade or other names:
25 (D2PM; diphenyl-2-pyrrolidinemethanol).

26 295. 3,4 Dichloromethylphenidate, some trade or
27 other names: (3,4-CTMP).

1 296. 3-chloromethyl-phenidate, some trade or other
2 names: (3-CTMP).
3 297. 4-Methylmethylphenidate.
4 298. 4-Fluoromethyl-phenidate, some trade or other
5 names: (4-FTMP).
6 299. Ethylphenidate.
7 300. Etizolam (Etilaam, Etizola, Sedekopan, Pasaden,
8 Depas).
9 301. Phenazepam.
10 302. Pyrazolam.
11 303. CL-218,872.
12 304. Zopiclone.
13 305. Salvinorin A.
14 306. AH-7921.
15 307. O-Desmethyltramadol, some trade or other names:
16 (O-DT; ODT).
17 308. Desmorphine (Dihydrodesoxymorphine; permonid;
18 krokodil; crocodile).
19 309. Acetyl Fentanyl (desmethylfentanyl).
20 310. 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine
21 (MT-45).
22 311. 1-(2-methoxyphenyl)piperazine, some trade or
23 other names: (MOPIP).
24 312. 1-(4-Chlorophenyl)piperazine, some trade or
25 other names: (pCPP).
26 313. para-Methoxyphenyl-piperazine, some trade or
27 other names: (MBZP).

1 314. Methylnmethaqualone.
2 315. Etaqualone.
3 316. 5-Iodo-2-aminoindane, some trade or other
4 names: (5-IAI).
5 317. 5,6-(Methylenedioxy)-2-aminoindane, some trade
6 or other names: (5,6-MDAI).
7 318. 4,5-(Methylenedioxy)-2-aminoindane, some trade
8 or other names: (4,5-MDAI).
9 319. MMAI.
10 320. W-15.
11 321. W-18.
12 322. Mitragynine.
13 323. Hydroxymitragynine.
14 324. Butyrfentanyl
15 (N-phenyl-N-[1-(2-phenylethyl)-4-pyridinyl]butyramide).
16 325. Beta-Hydroxythiofentanyl
17 (N-phenyl-N-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl}-pr
18 opanamide).
19 326. 4-methylphenethyl acetyl fentanyl
20 (N-phenyl-N-{1-[2-(4-methylphenyl)ethyl]-4-piperidinyl}-acetami
21 de).
22 327. Acrylfentanyl
23 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-prop-2-enamide).
24 328. 3-Allylfentanyl
25 (N-phenyl-N-[1-(2-phenylethyl)-(3s,4R)-3-prop-2-enyl-4-piperid
26 inyl]-propanamide).

1 329. Benzodioxole fentanyl
2 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-benzo[d][1,3]dioxole-5-carboxamide) .
3
4 330. Benzyl carfentanil
5 (N-phenyl-N-(1-benzyl-4-methylcarboxylate-4-piperidinyl)-propanamide) .
6
7 331. Brifentanil
8 (N-(2-fluorophenyl)-N-((3R,4S)-1-[2-(4-ethyl-5-oxotetrazol-1-yl)ethyl]-3-methyl-4-piperidinyl)-2-methoxyacetamide) .
9
10 332. Cyclopentylfentanyl
11 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-cyclopentanecarboxamide) .
12
13 333. 2,5-Dimethylfentanyl
14 (N-phenyl-N-[1-(2-phenylethyl)-2,5-dimethyl-4-piperidinyl]-propanamide) .
15
16 334. 4-Fluoroisobutyryl fentanyl
17 (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-isobutyramide) .
18
19 335. Furanyl fentanyl
20 (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]furan-2-carboxamide) .
21
22 336. Furanylethyl fentanyl
23 (N-phenyl-N-[1-(2-furanylethyl)-4-piperidinyl]-propanamide) .
24
25 337. Isobutyryl fentanyl
26 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-methylpropanamide) .

1 338. Lofentanil
2 (N-phenyl-N-[1-(2-phenylethyl)-(3R,4S)-3-methyl-4-methylcarbox
3 ylate-4-piperidinyl]-propanamide) .
4 339. 4-Methoxybutyrfentanyl
5 (N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butyr
6 amide) .
7 340. 4-Methoxymethylfentanyl
8 (N-phenyl-N-[1-(2-phenylethyl)-4-methoxymethyl-4-piperidinyl]-
9 propanamide) .
10 341. Meta-fluorobutyryl fentanyl
11 (N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-butyram
12 ide) .
13 342. Meta-fluorofentanyl
14 (N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-propa
15 mide) .
16 343. 3-Methylbutyrfentanyl
17 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butyramide) .
18 344. N-Methylcarfentanyl
19 (N-phenyl-N-(1-methyl-4-methylcarboxylate-4-piperidinyl)-propa
20 namide) .
21 345. Methoxyacetylfentanyl
22 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-methoxyacetami
23 de) .
24 346. Mirfentanyl
25 (N-(2-pyrazinyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-2-furamid
26 e) .

1 347. Ocfentanil
2 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperdiny]l)-2-metho
3 xyacetamide).

4 348. Ohmefentanil
5 (N-phenyl-N-[1-(2-hydroxy-2-phenylethyl)-3-methyl-4-piperidiny
6 l]-propanamide).

7 349. Ortho-fluorobutyryl fentanyl
8 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-butyram
9 ide).

10 350. Ortho-fluorofentanyl
11 (N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidnyl]-propa
12 mide).

13 351. Para-chlorofentanyl
14 (N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidiny]l)-propan
15 amide).

16 352. Para-chloroisobutyryl fentanyl
17 (N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidiny]l)-isobut
18 yramide).

19 353. 4-Fluorobutyryl fentanyl
20 (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidiny]l)-butyra
21 mide).

22 354. Para-methoxyfentanyl
23 (N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidiny]l)-propa
24 namide).

25 355. Para-methylfentanyl
26 (N-(4-methylphenyl)-N-[1-(2-phenylethyl)-4-piperidiny]l)-propan
27 amide).

1 356. 4-Phenyl fentanyl
2 (N-phenyl-N-[4-phenyl-1-(2-phenylethyl)-4-piperidinyl]-propana
3 mide) .
4 357. Trefentanyl
5 (N-(2-fluorophenyl)-N-{1-[2-(4-ethyl-5-oxo-4,5-dihydro-1H-tetr
6 azol-1-yl)ethyl]-4-phenyl-4-piperidinyl}-propanamide) .
7 358. Valeryl fentanyl
8 (N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide) .
9 359. Alpha-Methylacetylfentanyl
10 (N-phenyl-N-[1-phenylpropan-2-yl]-4-piperidinyl)-acetamide) .
11 360. Alpha-Methylbutyrfentanyl
12 (N-phenyl-N-[1-phenylpropan-2-yl]-4-piperidinyl)-butyramide) .
13 361. Alpha-Methylthiofentanyl
14 (N-phenyl-N-[1-(1-thienyl-2-ylpropan-2-yl)-4-piperidinyl]-prop
15 anamide) .
16 362. Beta-Hydroxy fentanyl
17 (N-phenyl-N-[1-(2-hydroxy-2-phenylethyl)-4-piperidinyl]-propan
18 amide) .
19 363. Beta-Methyl fentanyl
20 (N-phenyl-N-[1-(2-phenylpropyl)-4-piperidinyl]-propanamide) .
21 364. U-47700
22 (3,4-Dichloro-N-[(1R,2R)-2-(dimethylamino)cyclohexyl]-N-methyl
23 benzamide) .
24 365. W-19
25 ((Z)-N-{1-[2-(4-aminophenyl)ethyl]piperidin-2-ylidene}-4-chlor
26 obenzenesulfonamide) .

1 366. Flubromazolam
2 (8-bromo-6-(2-fluorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a]
3 [1,4]benzodiazepine).

4 377. Tianeptine.

5 (5)a. A synthetic controlled substance analogue,
6 being a material, mixture, or preparation that contains any
7 chemical structure of which is chemically similar to the
8 chemical structure of any other controlled substance in
9 Schedule I or Schedule II or that satisfies any one of the
10 following:

11 1. Has a stimulant, depressant, or hallucinogenic
12 effect on the central nervous system that mimics or is similar
13 to or greater than the stimulant, depressant, or
14 hallucinogenic effect on the central nervous system of a
15 controlled substance in Schedule I or Schedule II.

16 2. With respect to a particular person, if the
17 person represents or intends that the substance have a
18 stimulant, depressant, or hallucinogenic effect on the central
19 nervous system that is substantially similar to or greater
20 than the stimulant, depressant, or hallucinogenic effect on
21 the central nervous system of a controlled substance in
22 Schedule I or Schedule II and the substance is actually
23 capable of producing a stimulant, depressant, or
24 hallucinogenic effect on the central nervous system that
25 mimics, is similar to, or is greater than the stimulant,
26 depressant, or hallucinogenic effect on the central nervous
27 system of a controlled substance in Schedule I or Schedule II.

1 3. Has been demonstrated to have binding activity at
2 one or more cannabinoid receptors.

3 4. Is capable of exhibiting cannabinoid-like
4 activity.

5 5. Any compound structurally analogous to,
6 mimicking, or derived from 3-(1-naphthoyl)indole or
7 1H-indol-3-yl-(1-naphthyl)methane by substitution at the
8 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
9 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
10 cycloalkylmethyl, cycloalkylethyl,
11 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
12 or 2-(4-morpholinyl)ethyl whether or not further substituted
13 in the indole ring to any extent, whether or not substituted
14 in the naphthyl ring to any extent.

15 6. Any compound structurally analogous to,
16 mimicking, or derived from 3-(1-naphthoyl)pyrrole by
17 substitution at the nitrogen atom of the pyrrole ring by
18 alkyl, alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
19 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
20 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
21 or 2-(4-morpholinyl)ethyl, whether or not further substituted
22 in the pyrrole ring to any extent, whether or not substituted
23 in the naphthyl ring to any extent.

24 7. Any compound structurally analogous to,
25 mimicking, or derived from 1-(1-naphthylmethyl)indene by
26 substitution at the 3-position of the indene ring by alkyl,
27 alkyl halide, aryl halide, alkyl aryl halide, alkenyl,

1 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
2 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
3 or 2-(4-morpholinyl)ethyl whether or not further substituted
4 in the indene ring to any extent, whether or not substituted
5 in the naphthyl ring to any extent.

6 8. Any compound structurally analogous to,
7 mimicking, or derived from 3-phenylacetylindole by
8 substitution at the nitrogen atom of the indole ring with
9 alkyl, alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
10 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
11 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
12 or 2-(4-morpholinyl)ethyl, whether or not further substituted
13 in the indole ring to any extent, whether or not substituted
14 in the phenyl ring to any extent.

15 9. Any compound structurally analogous to,
16 mimicking, or derived from 2-(3-hydroxycyclohexyl)phenol by
17 substitution at the 5-position of the phenolic ring by alkyl,
18 alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
19 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
20 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
21 or 2-(4-morpholinyl)ethyl, whether or not substituted in the
22 cyclohexyl ring to any extent.

23 10. Any compound structurally analogous to,
24 mimicking, or derived from
25 3-(2,2,3,3-tetramethylcyclopropoyl)indole or
26 1H-indol-3-yl-(2,2,3,3-tetramethylcyclopropoyl)methane by
27 substitution at the nitrogen atom of the indole ring by alkyl,

1 alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
2 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
3 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
4 or 2-(4-morpholinyl)ethyl whether or not further substituted
5 in the indole ring to any extent.

6 11. Any compound structurally analogous to,
7 mimicking, or derived from 3-(adamant-1-oyl)indole or
8 1H-indol-3-yl-(1-adamantyl)methane by substitution at the
9 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
10 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
11 cycloalkylmethyl, cycloalkylethyl,
12 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
13 or 2-(4-morpholinyl)ethyl whether or not further substituted
14 in the indole ring to any extent.

15 12. Any compound structurally analogous to,
16 mimicking, or derived from
17 N-(1-naphthalenyl)indole-3-carboxamide or
18 1H-indol-(N-naphthyl)-3-carboxamide by substitution at the
19 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
20 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
21 cycloalkylmethyl, cycloalkylethyl,
22 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
23 or 2-(4-morpholinyl)ethyl whether or not further substituted
24 in the indole ring to any extent, whether or not substituted
25 in the naphthyl ring to any extent.

26 13. Any compound structurally analogous to,
27 mimicking, or derived from

1 N-(adamantan-1-yl)indole-3-carboxamide or
2 1H-indol-3-carboxamide-(1-adamantyl) by substitution at the
3 nitrogen atom of the indole ring by alkyl, alkyl halide, aryl
4 halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
5 cycloalkylmethyl, cycloalkylethyl,
6 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
7 or 2-(4-morpholinyl)ethyl whether or not further substituted
8 in the indole ring to any extent.

9 14. Any compound structurally analogous to,
10 mimicking, or derived from

11 N-(adamantan-1-yl)indazole-3-carboxamide or
12 1H-indazole-3-carboxamide-(1-adamantyl) by substitution at the
13 nitrogen atom of the indazole ring by alkyl, alkyl halide,
14 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
15 cycloalkylmethyl, cycloalkylethyl,
16 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
17 or 2-(4-morpholinyl)ethyl whether or not further substituted
18 in the indazole ring to any extent.

19 15. Any compound structurally analogous to,
20 mimicking, or derived from

21 N-[(1S)-1-(aminocarbonyl)-2-methylpropoyl]indazole-3-carboxyam
22 ide or
23 1H-indazole-3-carboxamide-N-[(1S)-1-(aminocarbonyl)-2-methylpr
24 opoyl] by substitution at the nitrogen atom of the indazole
25 ring by alkyl, alkyl halide, aryl halide, alkyl aryl halide,
26 alkenyl, aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
27 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,

1 or 2-(4-morpholinyl)ethyl whether or not further substituted
2 in the indazole ring to any extent.

3 16. Any compound structurally analogous to,
4 mimicking, or derived from 3-(1-naphthoyl)indazole or
5 1H-indazole-3-yl-(1-naphthyl)methane by substitution at the
6 nitrogen atom of the indazole ring by alkyl, alkyl halide,
7 aryl halide, alkyl aryl halide, alkenyl, aliphatic alcohol,
8 cycloalkylmethyl, cycloalkylethyl,
9 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
10 or 2-(4-morpholinyl)ethyl whether or not further substituted
11 in the indazole ring to any extent, whether or not substituted
12 in the naphthyl ring to any extent.

13 17. Any compound structurally analogous to,
14 mimicking, or derived from 3-(carboxylic acid 8-quinolinyl
15 ester)indole or 1H-indol-3-carboxylic acid-(8-quinolinyl)ester
16 by substitution at the nitrogen atom of the indole ring by
17 alkyl, alkyl halide, aryl halide, alkyl aryl halide, alkenyl,
18 aliphatic alcohol, cycloalkylmethyl, cycloalkylethyl,
19 (N-alkylpiperidin-2-yl)methyl, (tetrahydropyran-4-yl)methyl,
20 or 2-(4-morpholinyl)ethyl whether or not further substituted
21 in the indole ring to any extent, whether or not substituted
22 in the quinoline ring to any extent.

23 18. Any compound structurally related to
24 2-(4-iodo-2,5-dimethoxyphenyl) ethanamine by substitution of
25 the iodo moiety (4 position) with other halides, alkyl, alkyl
26 halides, thioalkyl, cycloalkyl, cycloalkylhalides and/or
27 substitution at the nitrogen atom of the ethanamine with

1 alkyl, alkyl halide, alkenyl, cycloalkylmethyl,
2 cycloalkylethyl, phenyl, benzyl whether or not further
3 substituted in the (either) phenyl ring to any extent.

4 19. Any compound structurally related to
5 2,5-dimethoxy-4-chloroamphetamine by substitution of the
6 chloro moiety (4 position) with other halides, alkyl, alkyl
7 halides, thioalkyl, cycloalkyl, cycloalkylhalides and/or
8 substitution at the nitrogen atom with alkyl, alkyl halide,
9 alkenyl, cycloalkylmethyl, cycloalkylethyl, phenyl, benzyl
10 whether or not further substituted in the (either) phenyl ring
11 to any extent.

12 20. Any compound structurally related to
13 2-amino-1-phenyl-1-propanone (cathinone) by substitution of
14 the amine with alkyl, alkyl halide, alkenyl, cycloalkylmethyl,
15 cycloalkylethyl, phenyl, benzyl whether or not further
16 substituted in the (either) phenyl ring to any extent.

17 21. Any compound structurally related to
18 a-pyrrolidinopentiophenone (a-pvp) whether or not further
19 substituted in the phenyl ring to any extent, whether or not
20 further substituted in the pyrrolidine ring to any extent.

21 b. A synthetic controlled substance or analogue in
22 subdivision (4) or this subdivision does not include any of
23 the following:

24 1. Any substance for which there is an approved new
25 drug application under the Federal Food, Drug, and Cosmetic
26 Act.

1 2. With respect to a particular person, any
2 substance, if an exemption is in effect for investigational
3 use, for that person, as provided by 21 U.S.C. § 355, and the
4 person is registered as a controlled substance researcher as
5 required under section 152.12, subdivision 3, to the extent
6 conduct with respect to the substance is pursuant to the
7 exemption and registration.

8 c. A controlled substance analogue is treated as a
9 controlled substance in Schedule I.

10 d. After the Alabama Department of Forensic Sciences
11 has determined a substance to be a synthetic controlled
12 substance analogue under this section, the department shall
13 notify the Alabama Department of Public Health with
14 information relevant to scheduling as provided by Section
15 20-2-20."