

Sixty-sixth  
Legislative Assembly  
of North Dakota

## ENGROSSED HOUSE BILL NO. 1113

Introduced by

Judiciary Committee

(At the request of the State Board of Pharmacy)

1 A BILL for an Act to amend and reenact subsection 18 of section 19-03.1-01, section  
2 19-03.1-05, subsection 7 of section 19-03.1-07, subsection 4 of section 19-03.1-09,  
3 subsection 7 of section 19-03.1-11, and subsection 5 of section 19-03.1-13 of the North Dakota  
4 Century Code, relating to the definition of marijuana and the scheduling of controlled  
5 substances; and to declare an emergency.

6 **BE IT ENACTED BY THE LEGISLATIVE ASSEMBLY OF NORTH DAKOTA:**

7 **SECTION 1. AMENDMENT.** Subsection 18 of section 19-03.1-01 of the North Dakota  
8 Century Code is amended and reenacted as follows:

9 18. "Marijuana" means all parts of the plant cannabis sativa L., whether growing or not;  
10 the seeds thereof; the ~~resinous product of the combustion~~ resin extracted from any part  
11 of the plant ~~cannabis~~; and every compound, manufacture, salt, derivative, mixture, or  
12 preparation of the plant ~~or~~ its seeds, or resin. The term does not include the mature  
13 stalks of the plant, fiber produced from the stalks, oil or cake made from the seeds of  
14 the plant, any other compound, manufacture, salt, derivative, mixture, or preparation of  
15 mature stalks, except the resin extracted therefrom, fiber, oil, or cake, or the sterilized  
16 seed of the plant which is incapable of germination. The term marijuana does not  
17 include hemp as defined in section 4.1-18-01.

18 **SECTION 2. AMENDMENT.** Section 19-03.1-05 of the North Dakota Century Code is  
19 amended and reenacted as follows:

20 **19-03.1-05. Schedule I.**

21 1. The controlled substances listed in this section are included in schedule I.  
22 2. Schedule I consists of the drugs and other substances, by whatever official name,  
23 common or usual name, chemical name, or brand name designated, listed in this  
24 section.

- 1           3. Opiates. Unless specifically excepted or unless listed in another schedule, any of the  
2           following opiates, including their isomers, esters, ethers, salts, and salts of isomers,  
3           esters, and ethers, whenever the existence of those isomers, esters, ethers, and salts  
4           is possible within the specific chemical designation:
- 5           a. Acetylmethadol.
  - 6           b. Allylprodine.
  - 7           c. Alphacetylmethadol.
  - 8           d. Alphameprodine.
  - 9           e. Alphamethadol.
  - 10          f. Benzethidine.
  - 11          g. Betacetylmethadol.
  - 12          h. Betameprodine.
  - 13          i. Betamethadol.
  - 14          j. Betaprodine.
  - 15          k. Clonitazene.
  - 16          l. Dextromoramide.
  - 17          m. Diampromide.
  - 18          n. Diethylthiambutene.
  - 19          o. Difenoxin.
  - 20          p. Dimenoxadol.
  - 21          q. Dimepheptanol.
  - 22          r. Dimethylthiambutene.
  - 23          s. Dioxaphetyl butyrate.
  - 24          t. Dipipanone.
  - 25          u. Ethylmethylthiambutene.
  - 26          v. Etonitazene.
  - 27          w. Etoxidine.
  - 28          x. Furethidine.
  - 29          y. Hydroxypethidine.
  - 30          z. Ketobemidone.
  - 31          aa. Levomoramide.

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- 1           bb. Levophenacymorphan.
- 2           cc. Morpheridine.
- 3           dd. MPPP (also known as 1-methyl-4-phenyl-4-propionoxypiperidine).
- 4           ee. Noracymethadol.
- 5           ff. Norlevorphanol.
- 6           gg. Normethadone.
- 7           hh. Norpipanone.
- 8           ii. PEPAP (1-(2-Phenylethyl)-4-Phenyl-4-acetoxypiperidine).
- 9           jj. Phenadoxone.
- 10          kk. Phenampromide.
- 11          ll. Phenomorphan.
- 12          mm. Phenoperidine.
- 13          nn. Piritramide.
- 14          oo. Proheptazine.
- 15          pp. Properidine.
- 16          qq. Propiram.
- 17          rr. Racemoramide.
- 18          ss. Tilidine.
- 19          tt. Trimeperidine.
- 20          uu. 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]cyclohexyl-N-methylbenzamide  
21             (also known as U-47700).
- 22          vv. 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine (also know as MT-45).
- 23          ww. 3,4-dichloro-N-[[1-(dimethylamino)cyclohexyl]methyl]benzamide (also known as  
24             AH-7921).
- 25          xx. Fentanyl derivatives. Unless specifically excepted or unless listed in another  
26             schedule or are not FDA approved drugs, and are derived from N-(1-(2-  
27             Phenylethyl)-4-piperidiny)-N-phenylpropanamide (Fentanyl) by any substitution  
28             on or replacement of the phenethyl group, any substitution on the piperidine ring,  
29             any substitution on or replacement of the propanamide group, any substitution on  
30             the anilido phenyl group, or any combination of the above. Examples include:

- 1 (1) N-[1-(1-methyl-2-phenethyl)-4-piperidiny]-N-phenylacetamide (also known  
2 as Acetyl-alpha-methylfentanyl).
- 3 (2) N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidy]propionanilide; 1-(1-methyl-  
4 2-phenylethyl)-4-(N-propanilido)piperidine (also known as Alpha-  
5 methylfentanyl).
- 6 (3) N-[1-methyl-2-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide (also  
7 known as Alpha-methylthiofentanyl).
- 8 (4) N-[1-(2-hydroxy-2-phenethyl)-4-piperidiny]-N-phenylpropanamide (also  
9 known as Beta-hydroxyfentanyl).
- 10 (5) N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidiny]-N-phenylpropanamide  
11 (also known as Beta-hydroxy-3-methylfentanyl).
- 12 (6) N-[3-methyl-1-(2-phenylethyl)-4-piperidy]-N-phenylpropanamide (also  
13 known as 3-Methylfentanyl).
- 14 (7) N-[3-methyl-1-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide (also  
15 known as 3-Methylthiofentanyl).
- 16 (8) N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidiny]propanamide (also  
17 known as Para-fluorofentanyl).
- 18 (9) N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidiny]propanamide (also known as  
19 Thiofentanyl).
- 20 (10) N-(1-phenylethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (also known  
21 as Furanyl Fentanyl).
- 22 (11) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide; N-(1-phenethylpiperidin-  
23 4-yl)-N-phenylbutanamide (also known as Butyryl Fentanyl).
- 24 (12) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide;  
25 N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidiny]-N-phenylpropanamide (also  
26 known as Beta-Hydroxythiofentanyl).
- 27 (13) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as Acetyl  
28 Fentanyl).
- 29 (14) N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]prop-2-enamideN-(1-phenethylpi  
30 peridin-4-yl)-N-phenylacrylamide (also known as AerylfentanylAcryl  
31 Fentanyl).

- 1 (15) N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]pentanamideN-(1-phenethylpip  
2 eridin-4-yl)-N-phenylpentanamide (also known as Valeryl Fentanyl).
- 3 (16) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known  
4 as 4-Fluoroisobutyryl Fentanyl).
- 5 (17) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide (also known  
6 as Ortho-fluorofentanyl, 2-Fluorofentanyl).
- 7 (18) N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (also  
8 known as Tetrahydrofuranyl Fentanyl).
- 9 (19) 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as  
10 Methoxyacetyl Fentanyl).
- 11 (20) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (also  
12 known as Cyclopropyl Fentanyl).
- 13 (21) N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide (also  
14 known as Ocfentanil).
- 15 (22) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (also  
16 known as Cyclopentyl Fentanyl).
- 17 (23) N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (also known as  
18 Isobutyryl Fentanyl).
- 19 (24) N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known  
20 as Para-chloroisobutyryl Fentanyl).
- 21 (25) N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known  
22 as Para-methoxybutyryl Fentanyl).
- 23 (26) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known as  
24 Para-fluorobutyryl Fentanyl).
- 25 4. Opium derivatives. Unless specifically excepted or unless listed in another schedule,  
26 any of the following opium derivatives, its salts, isomers, and salts of isomers  
27 whenever the existence of such salts, isomers, and salts of isomers is possible within  
28 the specific chemical designation:
- 29 a. Acetorphine.
- 30 b. Acetyldihydrocodeine.
- 31 c. Benzylmorphine.

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- 1 d. Codeine methylbromide.
- 2 e. Codeine-N-Oxide.
- 3 f. Cyrenorphine.
- 4 g. Desomorphine.
- 5 h. Dihydromorphine.
- 6 i. Drotebanol.
- 7 j. Etorphine (except hydrochloride salt).
- 8 k. Heroin.
- 9 l. Hydromorphinol.
- 10 m. Methyldesorphine.
- 11 n. Methyldihydromorphine.
- 12 o. Morphine methylbromide.
- 13 p. Morphine methylsulfonate.
- 14 q. Morphine-N-Oxide.
- 15 r. Myrophine.
- 16 s. Nicocodeine.
- 17 t. Nicomorphine.
- 18 u. Normorphine.
- 19 v. Pholcodine.
- 20 w. Thebacon.
- 21 5. Hallucinogenic substances. Unless specifically excepted or unless listed in another  
22 schedule, any material, compound, mixture, or preparation containing any quantity of  
23 the following hallucinogenic substances, including their salts, isomers, and salts of  
24 isomers whenever the existence of those salts, isomers, and salts of isomers is  
25 possible within the specific chemical designation (for purposes of this subsection only,  
26 the term "isomer" includes the optical, position, and geometric isomers):
  - 27 a. Alpha-ethyltryptamine, its optical isomers, salts, and salts of isomers (also known  
28 as etryptamine; a-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole).
  - 29 b. Alpha-methyltryptamine.
  - 30 c. 4-methoxyamphetamine (also known as 4-methoxy-a-methylphenethylamine;  
31 paramethoxyamphetamine; PMA).

- 1           d. N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-alpha-  
2           methyl-3,4(methylenedioxy)phenylamine, and N-hydroxy MDA.
- 3           e. Hashish.
- 4           f. Ibogaine (also known as 7-Ethyl-6, 6B, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-  
5           6, 9-methano-5 H-pyrido [1', 2':1,2] azepero (5,4-b) indole; Tabernanthe iboga).
- 6           g. Lysergic acid diethylamide.
- 7           h. Marijuana.
- 8           i. Parahexyl (also known as 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro- 6,6,9-trimethyl-  
9           6H-dibenzol[b,d]pyran; Synhexyl).
- 10          j. Peyote (all parts of the plant presently classified botanically as Lophophora  
11          williamsii Lemaire, whether growing or not, the seeds thereof, any extract from  
12          any part of such plant, and every compound, manufacture, salts, derivative,  
13          mixture, or preparation of such plant, its seeds, or its extracts).
- 14          k. N-ethyl-3-piperidyl benzilate.
- 15          l. N-methyl-3-piperidyl benzilate.
- 16          m. Psilocybin.
- 17          n. Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained in a  
18          plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of  
19          the substances contained in the cannabis plant, or in the resinous extractives of  
20          such plant, including synthetic substances, derivatives, and their isomers with  
21          similar chemical structure and pharmacological activity to those substances  
22          contained in the plant; excluding tetrahydrocannabinols found in hemp as defined  
23          by section 4.1-18-01; such as the following:
- 24           (1) Delta-1 cis or trans tetrahydrocannabinol, and their optical isomers. Other  
25           names: Delta-9-tetrahydrocannabinol.
- 26           (2) Delta-6 cis or trans tetrahydrocannabinol, and their optical isomers.
- 27           (3) Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers.
- 28          (Since nomenclature of these substances is not internationally standardized,  
29          compounds of these structures, regardless of numerical designation of atomic  
30          positions covered.)

1           o.   Cannabinoids, synthetic. It includes the chemicals and chemical groups listed  
2                   below, including their homologues, salts, isomers, and salts of isomers. The term  
3                   "isomer" includes the optical, position, and geometric isomers.

4           (1)   Indole carboxaldehydes. Any compound structurally derived from 1H-indole-  
5                   3-carboxaldehyde or 1H-2-carboxaldehyde substituted in both of the  
6                   following ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,  
7                   cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-  
8                   piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,  
9                   1-(N-methyl-3- morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo  
10                  benzyl group; and, at the hydrogen of the carboxaldehyde by a phenyl,  
11                  benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, pyrrolidinyl, piperazinyl, or  
12                  propionaldehyde group whether or not the compound is further modified to  
13                  any extent in the following ways:

14               (a)   Substitution to the indole ring to any extent; or

15               (b)   Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
16                   cyclopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group to any  
17                   extent; or

18               (c)   A nitrogen heterocyclic analog of the indole ring; or

19               (d)   A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
20                   adamantyl, or cyclopropyl ring.

21               (e)   Examples include:

22               [1]   1-Pentyl-3-(1-naphthoyl)indole - Other names: JWH-018 and  
23                   AM-678.

24               [2]   1-Butyl-3-(1-naphthoyl)indole - Other names: JWH-073.

25               [3]   1-Pentyl-3-(4-methoxy-1-naphthoyl)indole - Other names:  
26                   JWH-081.

27               [4]   1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole - Other names:  
28                   JWH-200.

29               [5]   1-Propyl-2-methyl-3-(1-naphthoyl)indole - Other names:  
30                   JWH-015.

31               [6]   1-Hexyl-3-(1-naphthoyl)indole - Other names: JWH-019.



- 1 [7] 1-Pentyl-3-(4-methyl-1-naphthoyl)indole - Other names:  
2 JWH-122.
- 3 [8] 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole - Other names: JWH-210.
- 4 [9] 1-Pentyl-3-(4-chloro-1-naphthoyl)indole - Other names:  
5 JWH-398.
- 6 [10] 1-(5-fluoropentyl)-3-(1-naphthoyl)indole - Other names:  
7 AM-2201.
- 8 [11] 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole - Other  
9 names: RCS-8.
- 10 [12] 1-Pentyl-3-(2-methoxyphenylacetyl)indole - Other names:  
11 JWH-250.
- 12 [13] 1-Pentyl-3-(2-methylphenylacetyl)indole - Other names:  
13 JWH-251.
- 14 [14] 1-Pentyl-3-(2-chlorophenylacetyl)indole - Other names: JWH-  
15 203.
- 16 [15] 1-Pentyl-3-(4-methoxybenzoyl)indole - Other names: RCS-4.
- 17 [16] (1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole) - Other names:  
18 AM-694.
- 19 [17] (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-  
20 yl]methanone - Other names: WIN 48,098 and Pravadoline.
- 21 [18] (1-Pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone --  
22 Other names: UR-144.
- 23 [19] (1-(5-fluoropentyl)indol-3-yl)-(2,2,3,3-  
24 tetramethylcyclopropyl)methanone - Other names: XLR-11.
- 25 [20] (1-(2-morpholin-4-ylethyl)-1H-indol-3-yl)-(2,2,3,3-  
26 tetramethylcyclopropyl)methanone - Other names: A-796,260.
- 27 [21] (1-(5-fluoropentyl)-1H-indazol-3-yl)(naphthalen-1-yl)methanone --  
28 Other names: THJ-2201.
- 29 [22] 1-naphthalenyl(1-pentyl-1H-indazol-3-yl)-methanone -- Other  
30 names: THJ-018.

- 1 [23] (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-yl)(naphthalen-1-  
2 yl)methanone - Other names: FUBIMINA.
- 3 [24] 1-[(N-methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl) indole -  
4 Other names: AM-1248.
- 5 [25] 1-Pentyl-3-(1-adamantoyl)indole - Other names: AB-001 and  
6 JWH-018 adamantyl analog.
- 7 (2) Indole carboxamides. Any compound structurally derived from 1H-indole-3-  
8 carboxamide or 1H-2-carboxamide substituted in both of the following ways:  
9 at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,  
10 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,  
11 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
12 morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo benzyl group;  
13 and, at the nitrogen of the carboxamide by a phenyl, benzyl, cumyl,  
14 naphthyl, adamantyl, cyclopropyl, or propionaldehyde group whether or not  
15 the compound is further modified to any extent in the following ways:
- 16 (a) Substitution to the indole ring to any extent; or
- 17 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
18 cyclopropyl, or propionaldehyde group to any extent; or
- 19 (c) A nitrogen heterocyclic analog of the indole ring; or
- 20 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
21 adamantyl, or cyclopropyl ring.
- 22 (e) Examples include:
- 23 [1] N-Adamantyl-1-pentyl-1H-indole-3-carboxamide - Other names:  
24 JWH-018 adamantyl carboxamide, APICA, SDB-001, and 2NE1.
- 25 [2] N-Adamantyl-1-fluoropentylindole-3-carboxamide - Other names:  
26 STS-135.
- 27 [3] N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide - Other  
28 names: AKB 48 and APINACA.
- 29 [4] N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide - Other  
30 names: NNEI and MN-24.

- 1 [5] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-  
2 carboxamide - Other names: ADBICA.
- 3 [6] (S)-N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-  
4 3-carboxamide - Other names: AB-PINACA.
- 5 [7] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-  
6 fluorophenyl)methyl]-1H-indazole-3-carboxamide - Other names:  
7 AB-FUBINACA.
- 8 [8] (~~S~~)-N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-  
9 indazole-3-carboxamide - Other names: 5-Fluoro AB-PINACA  
10 and 5F-AB-PINACA.
- 11 [9] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-  
12 3-carboxamide - Other names: ADB-PINACA.
- 13 [10] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl)-  
14 1H-indazole-3-carboxamide - Other names: AB-CHMINACA.
- 15 [11] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-  
16 indazole-3-carboxamide - Other names: ADB-FUBINACA.
- 17 [12] N-((3s,5s,7s)-adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-  
18 carboxamide - Other names: FUB-AKB48 and AKB48 N-(4-  
19 fluorobenzyl) analog.
- 20 [13] 1-(5-fluoropentyl)-N-(quinolin-8-yl)-1H-indazole-3-carboxamide -  
21 Other names: 5-fluoro-THJ.
- 22 [14] (~~S~~)-methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-  
23 methylbutanoate - Other names: 5-fluoro AMB and 5F-AMB.
- 24 [15] methyl (1-(4-fluorobenzyl)-1H-indazole-3-carbonyl)-L-valinate2-  
25 (1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-  
26 methylbutanoate - Other names: FUB-AMB, MMB-FUBINACA,  
27 and AMB-FUBINACA.
- 28 [16] N-[1-(aminocarbonyl)-2,2-dimethylpropyl]-1-(cyclohexylmethyl)-1  
29 H-indazole-3-carboxamide - Other names: MAB-CHMINACA and  
30 ADB-CHMINACA.

- 1 [17] Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-  
2 dimethylbutanoate - Other names: 5F-ADB and  
3 5F-MDMB-PINACA.
- 4 [18] N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-  
5 carboxamide - Other names: 5F-APINACA and 5F-AKB48.
- 6 [19] Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-  
7 dimethylbutanoate - Other names: MDMB-CHMICA and  
8 MMB-CHMINACA.
- 9 [20] Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-  
10 dimethylbutanoate - Other names: MDMB-FUBINACA.
- 11 [21] 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxa  
12 mide - Other names: 4-CN-CUMYL-BUTINACA; 4-cyano-  
13 CUMYL-BUTINACA; 4-CN-CUMYL BINACA; CUMYL-4CN-  
14 BINACA; SGT-78.
- 15 [22] methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-  
16 3-methylbutanoate - Other names: MMB-CHMICA, AMB-  
17 CHMICA.
- 18 [23] 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridi  
19 ne-3-carboxamide - Other names: 5F-CUMYL-P7AICA.
- 20 (3) Indole carboxylic acids. Any compound structurally derived from 1H-indole-  
21 3-carboxylic acid or 1H-2-carboxylic acid substituted in both of the following  
22 ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,  
23 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-  
24 piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,  
25 1-(N-methyl-3-morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo  
26 benzyl group; and, at the hydroxyl group of the carboxylic acid by a phenyl,  
27 benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, or propionaldehyde group  
28 whether or not the compound is further modified to any extent in the  
29 following ways:
- 30 (a) Substitution to the indole ring to any extent; or

- 1 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
2 cyclopropyl, propionaldehyde group to any extent; or  
3 (c) A nitrogen heterocyclic analog of the indole ring; or  
4 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
5 adamantyl, or cyclopropyl ring.  
6 (e) Examples include:  
7 [1] 1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid 8-quinolinyl  
8 ester - Other names: BB-22 and QUCHIC.  
9 [2] naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate -  
10 Other names: FDU-PB-22.  
11 [3] 1-pentyl-1H-indole-3-carboxylic acid 8-quinolinyl ester - Other  
12 names: PB-22 and QUPIC.  
13 [4] 1-(5-Fluoropentyl)-1H-indole-3-carboxylic acid 8-quinolinyl ester -  
14 Other names: 5-Fluoro PB-22 and 5F-PB-22.  
15 [5] quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate - Other  
16 names: FUB-PB-22.  
17 [6] naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate -  
18 Other names: NM2201 and CBL2201.  
19 (4) Naphthylmethylindoles. Any compound containing a 1H-indol-3-yl-(1-  
20 naphthyl)methane structure with substitution at the nitrogen atom of the  
21 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,  
22 cycloalkylethyl, 1-(N-methyl-2-piperidiny)methyl, 2-(4-morpholinyl)ethyl, 1-  
23 (N-methyl-2-pyrrolidiny)methyl, 1-(N-methyl-3-morpholinyl)methyl, or  
24 (tetrahydropyran-4-yl)methyl group whether or not further substituted in the  
25 indole ring to any extent and whether or not substituted in the naphthyl ring  
26 to any extent. Examples include:  
27 (a) 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane - Other names: JWH-175.  
28 (b) 1-Pentyl-1H-indol-3-yl-(4-methyl-1-naphthyl)methane - Other names:  
29 JWH-184.  
30 (5) Naphthoylpyrroles. Any compound containing a 3-(1-naphthoyl)pyrrole  
31 structure with substitution at the nitrogen atom of the pyrrole ring by an

- 1 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-  
2 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-  
3 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-  
4 yl)methyl group whether or not further substituted in the pyrrole ring to any  
5 extent, whether or not substituted in the naphthyl ring to any extent.  
6 Examples include: (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-  
7 ylmethanone - Other names: JWH-307.
- 8 (6) Naphthylmethylindenes. Any compound containing a naphthylideneindene  
9 structure with substitution at the 3-position of the indene ring by an alkyl,  
10 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-  
11 2-piperidinyl)methyl, 2 (4 morpholinyl)ethyl, 1-(N-methyl-2-  
12 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-  
13 yl)methyl group whether or not further substituted in the indene ring to any  
14 extent, whether or not substituted in the naphthyl ring to any extent.  
15 Examples include: E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane  
16 - Other names: JWH-176.
- 17 (7) Cyclohexylphenols. Any compound containing a 2-(3-  
18 hydroxycyclohexyl)phenol structure with substitution at the 5-position of the  
19 phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,  
20 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-  
21 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or  
22 (tetrahydropyran-4-yl)methyl group whether or not substituted in the  
23 cyclohexyl ring to any extent. Examples include:
- 24 (a) 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other  
25 names: CP 47,497.
- 26 (b) 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other  
27 names: Cannabicyclohexanol and CP 47,497 C8 homologue.
- 28 (c) 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-  
29 hydroxypropyl)cyclohexyl]-phenol - Other names: CP 55,940.
- 30 (8) Others specifically named:

- 1 (a) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-  
2 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names: HU-210.
- 3 (b) (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-  
4 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names:  
5 Dexanabinol and HU-211.
- 6 (c) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-  
7 benzoxazin-6-yl]-1-naphthalenylmethanone - Other names:  
8 WIN 55,212-2.
- 9 (d) Naphthalen-1-yl-(4-pentyloxynaphthalen-1-yl)methanone - Other  
10 names: CB-13.
- 11 p. Substituted phenethylamines. This includes any compound, unless specifically  
12 excepted, specifically named in this schedule, or listed under a different  
13 schedule, structurally derived from phenylethan-2-amine by substitution on the  
14 phenyl ring in any of the following ways, that is to say, by substitution with a fused  
15 methylenedioxy ring, fused furan ring, or fused tetrahydrofuran ring; by  
16 substitution with two alkoxy groups; by substitution with one alkoxy and either  
17 one fused furan, tetrahydrofuran, or tetrahydropyran ring system; or by  
18 substitution with two fused ring systems from any combination of the furan,  
19 tetrahydrofuran, or tetrahydropyran ring systems.
- 20 (1) Whether or not the compound is further modified in any of the following  
21 ways, that is to say:
- 22 (a) By substitution of phenyl ring by any halo, hydroxyl, alkyl,  
23 trifluoromethyl, alkoxy, or alkylthio groups;
- 24 (b) By substitution at the 2-position by any alkyl groups; or
- 25 (c) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl,  
26 hydroxybenzyl, methylenedioxybenzyl, or methoxybenzyl groups.
- 27 (2) Examples include:
- 28 (a) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (also known as 2C-C or  
29 2,5-Dimethoxy-4-chlorophenethylamine).
- 30 (b) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (also known as 2C-D or  
31 2,5-Dimethoxy-4-methylphenethylamine).

- 1 (c) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (also known as 2C-E or  
2 2,5-Dimethoxy-4-ethylphenethylamine).
- 3 (d) 2-(2,5-Dimethoxyphenyl)ethanamine (also known as 2C-H or 2,5-  
4 Dimethoxyphenethylamine).
- 5 (e) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (also known as 2C-I or  
6 2,5-Dimethoxy-4-iodophenethylamine).
- 7 (f) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (also known as 2C-N or  
8 2,5-Dimethoxy-4-nitrophenethylamine).
- 9 (g) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (also known as 2C-  
10 P or 2,5-Dimethoxy-4-propylphenethylamine).
- 11 (h) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (also known as 2C-  
12 T-2 or 2,5-Dimethoxy-4-ethylthiophenethylamine).
- 13 (i) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (also known as  
14 2C-T-4 or 2,5-Dimethoxy-4-isopropylthiophenethylamine).
- 15 (j) 2-(4-bromo-2,5-dimethoxyphenyl)ethanamine (also known as 2C-B or  
16 2,5-Dimethoxy-4-bromophenethylamine).
- 17 (k) 2-(2,5-dimethoxy-4-(methylthio)phenyl)ethanamine (also known as  
18 2C-T or 4-methylthio-2,5-dimethoxyphenethylamine).
- 19 (l) 1-(2,5-dimethoxy-4-iodophenyl)-propan-2-amine (also known as DOI  
20 or 2,5-Dimethoxy-4-iodoamphetamine).
- 21 (m) 1-(4-Bromo-2,5-dimethoxyphenyl)-2-aminopropane (also known as  
22 DOB or 2,5-Dimethoxy-4-bromoamphetamine).
- 23 (n) 1-(4-chloro-2,5-dimethoxy-phenyl)propan-2-amine (also known as  
24 DOC or 2,5-Dimethoxy-4-chloroamphetamine).
- 25 (o) 2-(4-bromo-2,5-dimethoxyphenyl)-N-[(2-  
26 methoxyphenyl)methyl]ethanamine (also known as 2C-B-NBOMe;  
27 2,5B-NBOMe or 2,5-Dimethoxy-4-bromo-N-(2-  
28 methoxybenzyl)phenethylamine).
- 29 (p) 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2 -  
30 methoxyphenyl)methyl]ethanamine (also known as 2C-I-NBOMe; 2,5I-



- 1 NBOMe or 2,5-Dimethoxy-4-iodo-N-(2-  
2 methoxybenzyl)phenethylamine).
- 3 (q) N-(2-Methoxybenzyl)-2-(3,4,5-trimethoxyphenyl)ethanamine (also  
4 known as mescaline-NBOMe or 3,4,5-trimethoxy-N-(2-  
5 methoxybenzyl)phenethylamine).
- 6 (r) 2-(4-chloro-2,5-dimethoxyphenyl)-N-[(2-  
7 methoxyphenyl)methyl]ethanamine (also known as 2C-C-NBOMe;  
8 2,5C-NBOMe or 2,5-Dimethoxy-4-chloro-N-(2-  
9 methoxybenzyl)phenethylamine).
- 10 (s) 2-(7-Bromo-5-methoxy-2,3-dihydro-1-benzofuran-4-yl)ethanamine  
11 (also known as 2CB-5-hemiFLY).
- 12 (t) 2-(8-bromo-2,3,6,7-tetrahydrofuro [2,3-f][1]benzofuran-4-  
13 yl)ethanamine (also known as 2C-B-FLY).
- 14 (u) 2-(10-Bromo-2,3,4,7,8,9-hexahydropyrano[2,3-g]chromen-5-  
15 yl)ethanamine (also known as 2C-B-butterFLY).
- 16 (v) N-(2-Methoxybenzyl)-1-(8-bromo-2,3,6,7-tetrahydrobenzo[1,2-b:4,5-  
17 b']difuran-4-yl)-2-aminoethane (also known as 2C-B-FLY-NBOMe).
- 18 (w) 1-(4-Bromofuro[2,3-f][1]benzofuran-8-yl)propan-2-amine (also known  
19 as bromo-benzodifuranyl-isopropylamine or bromo-dragonFLY).
- 20 (x) N-(2-Hydroxybenzyl)-4-iodo-2,5-dimethoxyphenethylamine (also  
21 known as 2C-I-NBOH or 2,5I-NBOH).
- 22 (y) 5-(2-Aminopropyl)benzofuran (also known as 5-APB).
- 23 (z) 6-(2-Aminopropyl)benzofuran (also known as 6-APB).
- 24 (aa) 5-(2-Aminopropyl)-2,3-dihydrobenzofuran (also known as 5-APDB).
- 25 (bb) 6-(2-Aminopropyl)-2,3,-dihydrobenzofuran (also known as 6-APDB).
- 26 (cc) 2,5-dimethoxy-amphetamine (also known as 2,5-dimethoxy-a-  
27 methylphenethylamine; 2,5-DMA).
- 28 (dd) 2,5-dimethoxy-4-ethylamphetamine (also known as DOET).
- 29 (ee) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (also known as 2C-T-  
30 7).
- 31 (ff) 5-methoxy-3,4-methylenedioxy-amphetamine.

- 1 (gg) 4-methyl-2,5-dimethoxy-amphetamine (also known as 4-methyl-2,5-  
2 dimethoxy- $\alpha$ -methylphenethylamine; DOM and STP).  
3 (hh) 3,4-methylenedioxy amphetamine (also known as MDA).  
4 (ii) 3,4-methylenedioxymethamphetamine (also known as MDMA).  
5 (jj) 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-  
6  $\alpha$ -methyl-3,4(methylenedioxy)phenethylamine, MDE, MDEA).  
7 (kk) 3,4,5-trimethoxy amphetamine.  
8 (ll) Mescaline (also known as 3,4,5-trimethoxyphenethylamine).
- 9 q. Substituted tryptamines. This includes any compound, unless specifically  
10 excepted, specifically named in this schedule, or listed under a different  
11 schedule, structurally derived from 2-(1H-indol-3-yl)ethanamine (i.e., tryptamine)  
12 by mono- or di-substitution of the amine nitrogen with alkyl or alkenyl groups or  
13 by inclusion of the amino nitrogen atom in a cyclic structure whether or not the  
14 compound is further substituted at the  $\alpha$ -position with an alkyl group or  
15 whether or not further substituted on the indole ring to any extent with any alkyl,  
16 alkoxy, halo, hydroxyl, or acetoxy groups. Examples include:
- 17 (1) 5-methoxy-N,N-diallyltryptamine (also known as 5-MeO-DALT).  
18 (2) 4-acetoxy-N,N-dimethyltryptamine (also known as 4-AcO-DMT or O-  
19 Acetylpsilocin).  
20 (3) 4-hydroxy-N-methyl-N-ethyltryptamine (also known as 4-HO-MET).  
21 (4) 4-hydroxy-N,N-diisopropyltryptamine (also known as 4-HO-DIPT).  
22 (5) 5-methoxy-N-methyl-N-isopropyltryptamine (also known as 5-MeO-MIPT).  
23 (6) 5-methoxy-N,N-dimethyltryptamine (also known as 5-MeO-DMT).  
24 (7) Bufotenine (also known as 3-(Beta-Dimethyl-aminoethyl)-5-hydroxyindole;  
25 3-(2-dimethylaminoethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-  
26 dimethyltryptamine; mappine).  
27 (8) 5-methoxy-N,N-diisopropyltryptamine (also known as 5-MeO-DiPT).  
28 (9) Diethyltryptamine (also known as N,N-Diethyltryptamine; DET).  
29 (10) Dimethyltryptamine (also known as DMT).  
30 (11) Psilocyn.
- 31 r. 1-[3-(trifluoromethylphenyl)]piperazine (also known as TFMPP).

- 1 s. 1-[4-(trifluoromethylphenyl)]piperazine.
- 2 t. 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (also known as 5,6-
- 3 Methylenedioxy-2-aminoindane or MDAI).
- 4 u. 2-(Ethylamino)-2-(3-methoxyphenyl)cyclohexanone (also known as
- 5 Methoxetamine or MXE).
- 6 v. Ethylamine analog of phencyclidine (also known as N-ethyl-1-
- 7 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl)
- 8 ethylamine, cyclohexamine, PCE).
- 9 w. Pyrrolidine analog of phencyclidine (also known as 1-(1-phenylcyclohexyl)-
- 10 pyrrolidine, PCPy, PHP).
- 11 x. Thiophene analog of phencyclidine (also known as (1-[1-(2-thienyl) cyclohexyl]
- 12 piperidine; 2-Thienylanalog of phencyclidine; TPCP, TCP).
- 13 y. 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (also known as TCPy).
- 14 z. Salvia divinorum, salvinorin A, or any of the active ingredients of salvia divinorum.
- 15 6. Depressants. Unless specifically excepted or unless listed in another schedule, any
- 16 material compound, mixture, or preparation which contains any quantity of the
- 17 following substances having a depressant effect on the central nervous system,
- 18 whenever the existence of such salts, isomers, and salts of isomers is possible within
- 19 the specific chemical designation:
  - 20 a. ~~Flunitrazepam.~~
  - 21 ~~b. Gamma-hydroxybutyric acid.~~
  - 22 ~~e.b. Mecloqualone.~~
  - 23 ~~d.c. Methaqualone.~~
- 24 7. Stimulants. Unless specifically excepted or unless listed in another schedule, any
- 25 material, compound, mixture, or preparation which contains any quantity of the
- 26 following substances having a stimulant effect on the central nervous system,
- 27 including its salts, isomers, and salts of isomers:
  - 28 a. Aminorex (also known as 2-amino-5-phenyl-2-oxazoline, or 4,5-dihydro-5-phenyl-
  - 29 2-oxazolamine).
  - 30 b. Cathinone.

1           c. Substituted cathinones. Any compound, material, mixture, preparation, or other  
2           product, unless listed in another schedule or an approved food and drug  
3           administration drug (e.g., bupropion, pyrovalerone), structurally derived from 2-  
4           aminopropan-1-one by substitution at the 1-position with either phenyl, naphthyl,  
5           or thiophene ring systems, whether or not the compound is further modified in  
6           any of the following ways:

- 7           (1) By substitution in the ring system to any extent with alkyl, alkylendioxy,  
8           alkoxy, haloalkyl, hydroxyl, or halide substituents, whether or not further  
9           substituted in the ring system by one or more other univalent substituents;  
10          (2) By substitution at the 3-position with an acyclic alkyl substituent;  
11          (3) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or  
12          methoxybenzyl groups; or  
13          (4) By inclusion of the 2-amino nitrogen atom in a cyclic structure.

14           Some trade or other names:

- 15          (a) 3,4-Methylenedioxy-alpha-pyrrolidinopropiophenone (also known as  
16           MDPPP).  
17          (b) 3,4-Methylenedioxy-N-ethylcathinone (also known as Ethylone,  
18           MDEC, or bk-MDEA).  
19          (c) 3,4-Methylenedioxy-N-methylcathinone (also known as Methylone or  
20           bk-MDMA).  
21          (d) 3,4-Methylenedioxy-pyrovalerone (also known as MDPV).  
22          (e) 3,4-Dimethylmethcathinone (also known as 3,4-DMMC).  
23          (f) 2-(methylamino)-1-phenylpentan-1-one (also known as Pentedrone).  
24          (g) 2-Fluoromethcathinone (also known as 2-FMC).  
25          (h) 3-Fluoromethcathinone (also known as 3-FMC).  
26          (i) 4-Methylethcathinone (also known as 4-MEC and 4-methyl-N-  
27           ethylcathinone).  
28          (j) 4-Fluoromethcathinone (also known as Flephedrone and 4-FMC).  
29          (k) 4-Methoxy-alpha-pyrrolidinopropiophenone (also known as MOPPP).  
30          (l) 4-Methoxymethcathinone (also known as Methedrone; bk-PMMA).  
31          (m) 4'-Methyl-alpha-pyrrolidinobutiophenone (also known as MPBP).

- 1 (n) Alpha-methylamino-butyrophenone (also known as Buphedrone or  
2 MABP).  
3 (o) Alpha-pyrrolidinobutiophenone (also known as alpha-PBP).  
4 (p) Alpha-pyrrolidinopropiophenone (also known as alpha-PPP).  
5 (q) Alpha-pyrrolidinopentiophenone (also known as Alpha-  
6 pyrrolidinovalerophenone or alpha-PVP).  
7 (r) Beta-keto-N-methylbenzodioxolylbutanamine (also known as Butylone  
8 or bk-MBDB).  
9 (s) Ethcathinone (also known as N-Ethylcathinone).  
10 (t) 4-Methylmethcathinone (also known as Mephedrone or 4-MMC).  
11 (u) Methcathinone.  
12 (v) N,N-dimethylcathinone (also known as metamfepramone).  
13 (w) Naphthylpyrovalerone (naphyrone).  
14 (x) B-Keto-Methylbenzodioxolylpentanamine (also known as Pentylone).  
15 (y) 4-Methyl-alpha-pyrrolidinopropiophenone (also known as 4-MePPP  
16 and MPPP).  
17 (z) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one (also known as  
18 Ephylone and N-Ethylpentylone).  
19 d. Fenethylline.  
20 e. Fluoroamphetamine.  
21 f. Fluoromethamphetamine.  
22 g. (±)cis-4-methylaminorex (also known as (±)cis-4,5-dihydro-4-methyl-5-phenyl-2-  
23 oxazolamine).  
24 h. N-Benzylpiperazine (also known as BZP, 1-benzylpiperazine).  
25 i. N-ethylamphetamine.  
26 j. N, N-dimethylamphetamine (also known as N,N-alpha-trimethyl-  
27 benzeneethanamine; N,N-alpha-trimethylphenethylamine).

28 **SECTION 3. AMENDMENT.** Subsection 7 of section 19-03.1-07 of the North Dakota  
29 Century Code is amended and reenacted as follows:

- 30 7. Hallucinogenic substances.

- 1           a. Nabilone [another name for nabilone ( $\pm$ )-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8,  
2                   10, 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9Hdibenzo [b, d] pyran-9-one].  
3           b. Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] in an oral solution in a drug  
4                   product approved for marketing by the federal food and drug administration.

5           **SECTION 4. AMENDMENT.** Subsection 4 of section 19-03.1-09 of the North Dakota  
6 Century Code is amended and reenacted as follows:

7           4. Depressants. Unless specifically excepted or unless listed in another schedule, any  
8                   material, compound, mixture, or preparation that contains any quantity of the following  
9                   substances having a depressant effect on the central nervous system:

10           a. Any compound, mixture, or preparation containing:

- 11                   (1) Amobarbital;  
12                   (2) Secobarbital;  
13                   (3) Pentobarbital;

14                   or any salt thereof and one or more other active medicinal ingredients which are  
15                   not listed in any schedule.

16           b. Any suppository dosage form containing:

- 17                   (1) Amobarbital;  
18                   (2) Secobarbital;  
19                   (3) Pentobarbital;

20                   or any salt of any of these drugs and approved by the food and drug  
21                   administration for marketing only as a suppository.

22           c. Any substance that contains any quantity of a derivative of barbituric acid, or any  
23                   salt of a derivative of barbituric acid, except those substances which are  
24                   specifically listed in other schedules thereof.

25           d. Chlorhexadol.

26           e. Embutramide.

27           f. Gamma-hydroxybutyric acid in a United States food and drug administration-  
28                   approved drug product.

29           g. Ketamine.

30           h. Lysergic acid.

31           i. Lysergic acid amide.

- 1           j. Methypylon.
- 2           k. Perampanel.
- 3           l. Sativex or its successor name as determined by the federal food and drug
- 4                 administration.
- 5           m. Sulfondiethylmethane.
- 6           ~~m.n.~~ Sulfonethylmethane.
- 7           ~~n.o.~~ Sulfonmethane.
- 8           ~~o.p.~~ Tiletamine and zolazepam or any salt thereof. Some trade or other names for a
- 9                 tiletamine-zolazepam combination product: Telazol. Some trade or other names
- 10                for tiletamine: 2-(ethylamino)-2-(2-thienyl)-cyclohexanone. Some trade or other
- 11                names for zolazepam: 4-2(2-fluorophenyl)-6, 8-dihydro-1,3,8-trimethylpyrazolo-
- 12                [3,4-e][1,4]-diazepin-7(1H)-one, flupyrazapon.

13           **SECTION 5. AMENDMENT.** Subsection 7 of section 19-03.1-11 of the North Dakota  
14 Century Code is amended and reenacted as follows:

- 15           7. Other substances. Unless specifically excepted or unless listed in another schedule,
- 16                 any material, compound, mixture, or preparation which contains any quantity of:
- 17                 a. Pentazocine, including its salts.
- 18                 b. Butorphanol, including its optical isomers.
- 19                 c. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-
- 20                 oxopropyl][(1S)-1-(4-phenyl-1*H*-imidazol-2-yl)ethyl]amino]methyl]-2-
- 21                 methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and
- 22                 salts of isomers.
- 23                 ~~d. Epidiolex or its successor name as determined by the United States food and~~
- 24                 ~~drug administration.~~

25           **SECTION 6. AMENDMENT.** Subsection 5 of section 19-03.1-13 of the North Dakota  
26 Century Code is amended and reenacted as follows:

- 27           5. Depressants. Unless specifically exempted or excluded or unless listed in another
- 28                 schedule, any material, compound, mixture, or preparation that contains any quantity
- 29                 of the following substances having a depressant effect on the central nervous system,
- 30                 including its salts, isomers, and salts of isomers whenever the existence of such salts,
- 31                 isomers, and salts of isomers is possible:

- 1           a. Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (also referred  
2           to as BRV; UCB-34714; Briviact) (including its salts).
- 3           b. Ezogabine N-[2-amino-4-(4-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester.
- 4           c. Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide].
- 5           d. Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid].
- 6           e. Approved cannabidiol drugs. A drug product in finished dosage formulation that  
7           has been approved by the federal food and drug administration, which contains  
8           cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-  
9           1,3-benzenediol) derived from cannabis and no more than 0.1 percent weight for  
10          weight residual tetrahydrocannabinols.
- 11          f. Gabapentin [2-[1-(aminomethyl) cyclohexyl] acetic acid].

12          **SECTION 7. EMERGENCY.** This Act is declared to be an emergency measure.