

**SENATE BILL NO. 2064  
with House Amendments**

**SENATE BILL NO. 2064**

Introduced by

Judiciary Committee

(At the request of the State Board of Pharmacy)

1 A BILL for an Act to amend and reenact sections 19-03.1-05, 19-03.1-09, and 19-03.1-11 of the  
2 North Dakota Century Code, relating to the scheduling of controlled substances; and to declare  
3 an emergency.

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

5 **SECTION 1. AMENDMENT.** Section 19-03.1-05 of the North Dakota Century Code is  
6 amended and reenacted as follows:

7 **19-03.1-05. Schedule I.**

- 8 1. The controlled substances listed in this section are included in schedule I.  
9 2. Schedule I consists of the drugs and other substances, by whatever official name,  
10 common or usual name, chemical name, or brand name designated, listed in this  
11 section.  
12 3. Opiates. Unless specifically excepted or unless listed in another schedule, any of the  
13 following opiates, including their isomers, esters, ethers, salts, and salts of isomers,  
14 esters, and ethers, whenever the existence of those isomers, esters, ethers, and salts  
15 is possible within the specific chemical designation:  
16 a. Acetylmethadol.  
17 b. Allylprodine.  
18 c. Alphacetylmethadol.  
19 d. Alphameprodine.  
20 e. Alphamethadol.  
21 f. Benzethidine.  
22 g. Betacetylmethadol.

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- 1 h. Betameprodine.
- 2 i. Betamethadol.
- 3 j. Betaprodine.
- 4 k. Brorphine.
- 5 l. Clonitazene.
- 6 m. Dextromoramide.
- 7 n. Diampromide.
- 8 o. Diethylthiambutene.
- 9 p. Difenoxin.
- 10 q. Dimenoxadol.
- 11 r. Dimepheptanol.
- 12 s. Dimethylthiambutene.
- 13 t. Dioxaphetyl butyrate.
- 14 u. Dipipanone.
- 15 v. Ethylmethylthiambutene.
- 16 w. Etonitazene.
- 17 x. Etoxidine.
- 18 y. Furethidine.
- 19 z. Hydroxypethidine.
- 20 aa. Isotonitazene (also known as N,N-diethyl-2-(2-(4- isopropoxybenzyl)-5-nitro-1H-
- 21 benzimidazol-1-yl)ethan-1-amine).
- 22 bb. Ketobemidone.
- 23 cc. Levomoramide.
- 24 dd. Levophenacymorphan.
- 25 ee. Morpheridine.
- 26 ff. MPPP (also known as 1-methyl-4-phenyl-4-propionoxypiperidine).
- 27 gg. Noracymethadol.
- 28 hh. Norlevorphanol.
- 29 ii. Normethadone.
- 30 jj. Norpipanone.
- 31 kk. PEPAP (1-(2-Phenylethyl)-4-Phenyl-4-acetoxypiperidine).

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- 1           ll.   Phenadoxone.
- 2           mm.   Phenampromide.
- 3           nn.   Phenomorphane.
- 4           oo.   Phenoperidine.
- 5           pp.   Piritramide.
- 6           qq.   Proheptazine.
- 7           rr.   Properidine.
- 8           ss.   Propiram.
- 9           tt.   Racemoramide.
- 10          uu.   Tilidine.
- 11          vv.   Trimeperidine.
- 12          ww.   3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as
- 13                    U-47700).
- 14          xx.   1-cyclohexyl-4-(1,2-diphenylethyl)piperazine (also known as MT-45).
- 15          yy.   3,4-dichloro-N-[[1-(dimethylamino)cyclohexyl]methyl]benzamide (also known as
- 16                    AH-7921).
- 17          zz.   Zipeprol.
- 18          aaa.   2-(2-(4-butoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine
- 19                    (also known as Butonitazene).
- 20          bbb.   2-(2-(4-ethoxybenzyl)-1H-benzimidazol-1-yl)-N,N-diethylethan-1-amine (also
- 21                    known as Etodesnitazene and etazene).
- 22          ccc.   N,N-diethyl-2-(2-(4-fluorobenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine
- 23                    (also known as Flunitazene).
- 24          ddd.   N,N-diethyl-2-(2-(4-methoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine (also
- 25                    known as Metodesnitazene).
- 26          eee.   N,N-diethyl-2-(2-(4-methoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine
- 27                    (also known as Metonitazene).
- 28          fff.   2-(4-ethoxybenzyl)-5-nitro-1-(2-(pyrrolidin-1-yl)ethyl)-1H-benzimidazole (also
- 29                    known as N-Pyrrolidino Etonitazene and Etonitazepyne).
- 30          ggg.   N,N-diethyl-2-(5-nitro-2-(4-propoxybenzyl)-1H-benzimidazol-1-yl)ethan-1-amine
- 31                    (also known as Protonitazene).



- 1 (10) N-(1-phenylethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (also known  
2 as Furanyl Fentanyl).
- 3 (11) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide; N-(1-phenethylpiperidin-  
4 4-yl)-N-phenylbutanamide (also known as Butyryl Fentanyl).
- 5 (12) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide;  
6 N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl]-N-phenylpropanamide (also  
7 known as Beta-Hydroxythiofentanyl).
- 8 (13) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as Acetyl  
9 Fentanyl).
- 10 (14) N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (also known as Acryl  
11 Fentanyl).
- 12 (15) N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (also known as Valeryl  
13 Fentanyl).
- 14 (16) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known  
15 as 4-Fluoroisobutyryl Fentanyl).
- 16 (17) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide (also known  
17 as Ortho-fluorofentanyl, 2-Fluorofentanyl).
- 18 (18) N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (also  
19 known as Tetrahydrofuran Fentanyl).
- 20 (19) 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as  
21 Methoxyacetyl Fentanyl).
- 22 (20) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (also  
23 known as Cyclopropyl Fentanyl).
- 24 (21) N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide (also  
25 known as Ocfentanil).
- 26 (22) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (also  
27 known as Cyclopentyl Fentanyl).
- 28 (23) N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (also known as  
29 Isobutyryl Fentanyl).
- 30 (24) N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known  
31 as Para-chloroisobutyryl Fentanyl).

- 1 (25) N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known  
2 as Para-methoxybutyryl Fentanyl).
- 3 (26) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known as  
4 Para-fluorobutyryl Fentanyl).
- 5 (27) N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide (also  
6 known as 2'-fluoro Ortho-fluorofentanyl; 2'-fluoro 2-fluorofentanyl).
- 7 (28) N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide (also known as  
8 Ortho-methyl Acetylfentanyl; 2-methyl acetylfentanyl).
- 9 (29) N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide (also known as  
10 Beta'-phenyl Fentanyl; 3-phenylpropanoyl fentanyl and Hydrocinnamoyl  
11 Fentanyl).
- 12 (30) N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide (also  
13 known as Thiofuranyl Fentanyl; 2-thiofuranyl fentanyl; thiophene fentanyl).
- 14 (31) (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide (also known as  
15 Crotonyl Fentanyl).
- 16 (32) N-(1-(4-methylphenethyl)piperidin-4-yl)-N-phenylacetamide (4'-methyl acetyl  
17 fentanyl).
- 18 (33) N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-yl)propionamide (beta-methyl  
19 fentanyl).
- 20 (34) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (ortho-  
21 fluorobutyryl fentanyl; 2-fluorobutyryl fentanyl).
- 22 (35) 2-methoxy-N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide  
23 (ortho-methyl methoxyacetylfentanyl; 2-methyl methoxyacetyl fentanyl).
- 24 (36) N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-yl)propionamide (para-  
25 methylfentanyl; 4-methylfentanyl).
- 26 (37) N-(1-phenethylpiperidin-4-yl)-N-phenylbenzamide (phenyl fentanyl; benzoyl  
27 fentanyl).
- 28 (38) Ethyl (1-phenethylpiperidin-4-yl)(phenyl)carbamate (fentanyl carbamate).
- 29 (39) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)acrylamide (ortho-fluoroacryl  
30 fentanyl).

- 1 (40) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (ortho-  
2 fluoroisobutyryl fentanyl).
- 3 (41) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide (para-  
4 fluoro furanyl fentanyl).
- 5 (42) 2',5'-dimethoxyfentanyl(N-(1-(2,5-dimethoxyphenethyl)piperidine-4-yl)-N-  
6 phenylpropionamide).
- 7 (43) 3-furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3-  
8 carboxamide).
- 9 (44) alpha'-methyl butyryl fentanyl(2-methyl-N-(1-phenethylpiperidin-4-yl)-N-  
10 phenylbutanamide).
- 11 (45) beta-methylacetyl fentanyl(N-phenyl-N-(1-(2-phenylpropyl)piperidin-4-  
12 yl)acetamide).
- 13 (46) isovaleryl fentanyl(3-methyl-N-(1-phenethylpiperidin-4-yl)-N-  
14 phenylbutanamide).
- 15 (47) meta-fluorofentanyl(N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-  
16 yl)propionamide).
- 17 (48) meta-fluorofuranyl fentanyl(N-3-fluorophenyl)-N-(1-phenethylpiperidin-4-  
18 yl)furan-2-carboxamide).
- 19 (49) meta-fluoroisobutyryl fentanyl(N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-  
20 yl)isobutyramide).
- 21 (50) ortho-chlorofentanyl(N-(2-chlorophenyl)-N-(1-phenethylpiperidin-4-  
22 yl)propionamide).
- 23 (51) ortho-fluorofuranyl fentanyl(N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-  
24 yl)furan-2-carboxamide).
- 25 (52) ortho-methylcyclopropylfentanyl(N-2-methylphenyl)-N-(1-phenethylpiperidin-  
26 4-yl)cyclopropanecarboxamide).
- 27 (53) para-chlorofentanyl(N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-  
28 yl)propionamide).
- 29 (54) para-fluoro valeryl fentanyl(N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-  
30 yl)pentanamide).

1                   (55) para-methoxyfuranyl fentanyl(N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-  
2   4-yl)furan-2-carboxamide).

3                   (56) para-methylcyclopropyl fentanyl(N-(4-methylphenyl)-N-(1-phenylpiperidin-4-  
4   yl)cyclopropanecarboxamide).

5                   (57) tetrahydrothiofuranyl fentanyl(N-(1-phenethylpiperidin-4-yl)-N-  
6   phenyltetrahydrothiophene-2-carboxamide).

7           4.   Opium derivatives. Unless specifically excepted or unless listed in another schedule,  
8                   any of the following opium derivatives, its salts, isomers, and salts of isomers  
9                   whenever the existence of such salts, isomers, and salts of isomers is possible within  
10                   the specific chemical designation:

- 11           a.   Acetorphine.
- 12           b.   Acetyldihydrocodeine.
- 13           c.   Benzylmorphine.
- 14           d.   Codeine methylbromide.
- 15           e.   Codeine-N-Oxide.
- 16           f.   Cyprenorphine.
- 17           g.   Desomorphine.
- 18           h.   Dihydromorphine.
- 19           i.   Drotebanol.
- 20           j.   Etorphine (except hydrochloride salt).
- 21           k.   Heroin.
- 22           l.   Hydromorphanol.
- 23           m.   Methyldesorphine.
- 24           n.   Methyldihydromorphine.
- 25           o.   Morphine methylbromide.
- 26           p.   Morphine methylsulfonate.
- 27           q.   Morphine-N-Oxide.
- 28           r.   Myrophine.
- 29           s.   Nicocodeine.
- 30           t.   Nicomorphine.
- 31           u.   Normorphine.



- 1 v. Pholcodine.
- 2 w. Thebacon.
- 3 5. Hallucinogenic substances. Unless specifically excepted or unless listed in another  
4 schedule, any material, compound, mixture, or preparation containing any quantity of  
5 the following hallucinogenic substances, including their salts, isomers, and salts of  
6 isomers whenever the existence of those salts, isomers, and salts of isomers is  
7 possible within the specific chemical designation (for purposes of this subsection only,  
8 the term "isomer" includes the optical, position, and geometric isomers):
- 9 a. Alpha-ethyltryptamine, its optical isomers, salts, and salts of isomers (also known  
10 as etryptamine;  $\alpha$ -ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole).
- 11 b. Alpha-methyltryptamine.
- 12 c. 4-methoxyamphetamine (also known as 4-methoxy- $\alpha$ -methylphenethylamine;  
13 paramethoxyamphetamine; PMA).
- 14 d. N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy- $\alpha$ -  
15 methyl-3,4(methylenedioxy)phenylamine, and N-hydroxy MDA).
- 16 e. Ibogaine (also known as 7-Ethyl-6, 6B, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-  
17 6, 9-methano-5 H-pyrido [1', 2':1,2] azepino (5,4-b) indole; Tabernanthe iboga).
- 18 f. Lysergic acid diethylamide.
- 19 g. Marijuana.
- 20 h. Parahexyl (also known as 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro- 6,6,9-trimethyl-  
21 6H-dibenzol[b,d]pyran; Synhexyl).
- 22 i. Peyote (all parts of the plant presently classified botanically as *Lophophora*  
23 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from  
24 any part of such plant, and every compound, manufacture, salts, derivative,  
25 mixture, or preparation of such plant, its seeds, or its extracts).
- 26 j. N-ethyl-3-piperidyl benzilate.
- 27 k. N-methyl-3-piperidyl benzilate.
- 28 l. Psilocybin.
- 29 m. (1) Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained  
30 in a plant of the genus *Cannabis* (cannabis plant), as well as synthetic  
31 equivalents of the substances contained in the cannabis plant, or in the

1 resinous extractives of such plant, including synthetic substances,  
2 derivatives, and their isomers with similar chemical structure and  
3 pharmacological activity to those substances contained in the plant; such as  
4 the following:

5 (a) Delta-1 cis or trans tetrahydrocannabinol, and their optical isomers.

6 Other names: Delta-9-tetrahydrocannabinol.

7 (b) Delta-6 cis or trans tetrahydrocannabinol, and their optical isomers.

8 Other names: Delta-8-tetrahydrocannabinol.

9 (c) Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers.

10 (Since nomenclature of these substances is not internationally standardized,  
11 compounds of these structures, regardless of numerical designation of atomic  
12 positions covered.)

13 (2) Tetrahydrocannabinols do not include:

14 (a) The allowable amount of total tetrahydrocannabinol found in hemp or  
15 an allowed hemp commodity or product as defined in chapter  
16 4.1-18.1; or

17 (b) A prescription drug approved by the United States food and drug  
18 administration under section 505 of the Federal Food, Drug, and  
19 Cosmetic Act [21 U.S.C. 355].

20 n. Cannabinoids, synthetic. It includes the chemicals and chemical groups listed  
21 below, including their homologues, salts, isomers, and salts of isomers. The term  
22 "isomer" includes the optical, position, and geometric isomers.

23 (1) Indole acetamides. Any compound structurally derived from 1H-indole3-  
24 acetamide or 1H-2-acetamide substituted in both of the following ways: at  
25 the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,  
26 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2piperidiny)methyl, 2-  
27 (4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidiny)methyl, 1-(N-methyl-3-  
28 morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo benzyl group;  
29 and, at the hydrogen of the acetamide by a phenyl, benzyl, cumyl, naphthyl,  
30 adamantyl, cyclopropyl, pyrrolidiny, piperaziny, or propionaldehyde group

- 1                   whether or not the compound is further modified to any extent in the  
2                   following ways:
- 3                   (a)    Substitution to the indole ring to any extent; or  
4                   (b)    Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
5                   cyclopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group to any  
6                   extent; or  
7                   (c)    A nitrogen heterocyclic analog of the indole ring; or  
8                   (d)    A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
9                   adamantyl, or cyclopropyl ring.  
10                  (e)    Examples include:
- 11                   [1]    N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide - Other names: CH-  
12                   PIATA, Cyclohexyl-PIATA, CHX-PIATA, CH-PIACA, and CHX-  
13                   PIACA.  
14                   [2]    N-cyclohexyl-2-[1-[(4-fluorophenyl)methyl]indol-3-yl]acetamide -  
15                   Other names: CH-FUBIATA and CH-FUBIACA.  
16                   [3]    2-[[2-[1-[(4-fluorophenyl)methyl]indol-3-yl]acetyl]amino]-3,3-  
17                   dimethyl-butanamide - Other names: ADB-FUBIATA, FUB-  
18                   ACADB, and AD-18.
- 19                  (2)    Indole carboxaldehydes. Any compound structurally derived from 1H-indole-  
20                  3-carboxaldehyde or 1H-2-carboxaldehyde substituted in both of the  
21                  following ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,  
22                  cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-  
23                  piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,  
24                  1-(N-methyl-3-morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo  
25                  benzyl group; and, at the hydrogen of the carboxaldehyde by a phenyl,  
26                  benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, pyrrolidinyl, piperazinyl, or  
27                  propionaldehyde group whether or not the compound is further modified to  
28                  any extent in the following ways:  
29                  (a)    Substitution to the indole ring to any extent; or

- 1 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
2 cyclopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group to any  
3 extent; or
- 4 (c) A nitrogen heterocyclic analog of the indole ring; or
- 5 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
6 adamantyl, or cyclopropyl ring.
- 7 (e) Examples include:
- 8 [1] 1-Pentyl-3-(1-naphthoyl)indole - Other names: JWH-018 and  
9 AM-678.
- 10 [2] 1-Butyl-3-(1-naphthoyl)indole - Other names: JWH-073.
- 11 [3] 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole - Other names:  
12 JWH-081.
- 13 [4] 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole - Other names:  
14 JWH-200.
- 15 [5] 1-Propyl-2-methyl-3-(1-naphthoyl)indole - Other names:  
16 JWH-015.
- 17 [6] 1-Hexyl-3-(1-naphthoyl)indole - Other names: JWH-019.
- 18 [7] 1-Pentyl-3-(4-methyl-1-naphthoyl)indole - Other names:  
19 JWH-122.
- 20 [8] 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole - Other names: JWH-210.
- 21 [9] 1-Pentyl-3-(4-chloro-1-naphthoyl)indole - Other names:  
22 JWH-398.
- 23 [10] 1-(5-fluoropentyl)-3-(1-naphthoyl)indole - Other names:  
24 AM-2201.
- 25 [11] 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole - Other  
26 names: RCS-8.
- 27 [12] 1-Pentyl-3-(2-methoxyphenylacetyl)indole - Other names:  
28 JWH-250.
- 29 [13] 1-Pentyl-3-(2-methylphenylacetyl)indole - Other names:  
30 JWH-251.

- 1 [14] 1-Pentyl-3-(2-chlorophenylacetyl)indole - Other names: JWH-  
2 203.
- 3 [15] 1-Pentyl-3-(4-methoxybenzoyl)indole - Other names: RCS-4.
- 4 [16] (1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole) - Other names:  
5 AM-694.
- 6 [17] (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-  
7 yl]methanone - Other names: WIN 48,098 and Pravadoline.
- 8 [18] (1-Pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone --  
9 Other names: UR-144.
- 10 [19] (1-(5-fluoropentyl)indol-3-yl)-(2,2,3,3-  
11 tetramethylcyclopropyl)methanone - Other names: XLR-11.
- 12 [20] (1-(2-morpholin-4-ylethyl)-1H-indol-3-yl)-(2,2,3,3-  
13 tetramethylcyclopropyl)methanone - Other names: A-796,260.
- 14 [21] (1-(5-fluoropentyl)-1H-indazol-3-yl)(naphthalen-1-yl)methanone --  
15 Other names: THJ-2201.
- 16 [22] 1-naphthalenyl(1-pentyl-1H-indazol-3-yl)-methanone -- Other  
17 names: THJ-018.
- 18 [23] (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-yl)(naphthalen-1-  
19 yl)methanone - Other names: FUBIMINA.
- 20 [24] 1-[(N-methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl) indole -  
21 Other names: AM-1248.
- 22 [25] 1-Pentyl-3-(1-adamantoyl)indole - Other names: AB-001 and  
23 JWH-018 adamantyl analog.
- 24 (3) Indole carboxamides. Any compound structurally derived from 1H-indole-3-  
25 carboxamide or 1H-2-carboxamide substituted in both of the following ways:  
26 at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,  
27 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,  
28 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-  
29 morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo benzyl group;  
30 and, at the nitrogen of the carboxamide by a phenyl, benzyl, cumyl,

- 1 naphthyl, adamantyl, cyclopropyl, or propionaldehyde group whether or not  
2 the compound is further modified to any extent in the following ways:
- 3 (a) Substitution to the indole ring to any extent; or  
4 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
5 cyclopropyl, or propionaldehyde group to any extent; or  
6 (c) A nitrogen heterocyclic analog of the indole ring; or  
7 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
8 adamantyl, or cyclopropyl ring.  
9 (e) Examples include:
- 10 [1] N-Adamantyl-1-pentyl-1H-indole-3-carboxamide - Other names:  
11 JWH-018 adamantyl carboxamide, APICA, SDB-001, and 2NE1.  
12 [2] N-Adamantyl-1-fluoropentylindole-3-carboxamide - Other names:  
13 STS-135.  
14 [3] N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide - Other  
15 names: AKB 48 and APINACA.  
16 [4] N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide - Other  
17 names: NNEI and MN-24.  
18 [5] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-  
19 carboxamide - Other names: ADBICA.  
20 [6] (S)-N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-  
21 3-carboxamide - Other names: AB-PINACA.  
22 [7] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-  
23 fluorophenyl)methyl]-1H-indazole-3-carboxamide - Other names:  
24 AB-FUBINACA.  
25 [8] N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-  
26 indazole-3-carboxamide - Other names: 5-Fluoro AB-PINACA  
27 and 5F-AB-PINACA.  
28 [9] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-  
29 3-carboxamide - Other names: ADB-PINACA.  
30 [10] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl)-  
31 1H-indazole-3-carboxamide - Other names: AB-CHMINACA.

- 1 [11] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-  
2 indazole-3-carboxamide - Other names: ADB-FUBINACA.
- 3 [12] N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H -indazole-3-  
4 carboxamide - Other names: FUB-AKB48, FUB-APINACA, and  
5 AKB48 N-(4-FLUOROBENZYL).
- 6 [13] 1-(5-fluoropentyl)-N-(quinolin-8-yl)-1H-indazole-3-carboxamide -  
7 Other names: 5-fluoro-THJ.
- 8 [14] methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-  
9 methylbutanoate - Other names: 5-fluoro AMB and 5F-AMB.
- 10 [15] methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-  
11 methylbutanoate - Other names: FUB-AMB, MMB-FUBINACA,  
12 and AMB-FUBINACA.
- 13 [16] N-[1-(aminocarbonyl)-2,2-dimethylpropyl]-1-(cyclohexylmethyl)-1  
14 H-indazole-3-carboxamide - Other names: MAB-CHMINACA and  
15 ADB-CHMINACA.
- 16 [17] Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-  
17 dimethylbutanoate - Other names: 5F-ADB and  
18 5F-MDMB-PINACA.
- 19 [18] N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-  
20 carboxamide - Other names: 5F-APINACA and 5F-AKB48.
- 21 [19] Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-  
22 dimethylbutanoate - Other names: MDMB-CHMICA and  
23 MMB-CHMINACA.
- 24 [20] Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-  
25 dimethylbutanoate - Other names: MDMB-FUBINACA.
- 26 [21] 1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxa  
27 mide - Other names: 4-CN-CUMYL-BUTINACA; 4-cyano-  
28 CUMYL-BUTINACA; 4-CN-CUMYL BINACA; CUMYL-4CN  
29 -BINACA; SGT-78.

- 1 [22] methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-  
2 3-methylbutanoate - Other names: MMB-CHMICA, AMB-  
3 CHMICA.
- 4 [23] 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyridi  
5 ne-3-carboxamide - Other names: 5F-CUMYL-P7AICA.
- 6 [24] ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-  
7 dimethylbutanoate - Other names: 5F-EDMB-PINACA.
- 8 [25] methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-  
9 dimethylbutanoate - Other names: 5F-MDMB-PICA and 5F-  
10 MDMB-2201.
- 11 [26] 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-  
12 carboxamide - Other names: 5F-CUMYL-PINACA, SGT-25.
- 13 [27] (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)  
14 methanone - Other names: FUB-144.
- 15 [28] methyl 2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-  
16 dimethylbutanoate (4F-MDMB-BINACA, 4F-MDMB-BUTINACA).
- 17 [29] ~~Methyl 3,3-dimethyl-2-[(1-pent-4-enylindazole-3-~~  
18 ~~carbonyl)amino]butanoate - Other names: MDMB-4en-PINACA,~~  
19 ~~MDMB-PENINACA, and 5-CL-ADB-A.~~
- 20 [30] Methyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-  
21 dimethyl-butanoate - Other names: 5F-MDMB-PICA and 5F-  
22 MDMB-2201.
- 23 [31] ~~1-butyl-N-(1-carbamoyl-2,2-dimethyl-propyl)indazole-3-~~  
24 ~~carboxamide - Other names: ADB-BINACA and ADB-BUTINACA.~~
- 25 [32][30] 5-bromo-N-(1-carbamoyl-2,2-dimethyl-propyl)-1H-indazole-3-  
26 carboxamide - Other names: ADB-5Br-INACA.
- 27 [33][31] Methyl 2-[(5-bromo-1H-indazole-3-carbonyl)amino]-3,3-dimethyl-  
28 butanoate - Other names: MDMB-5Br-INACA.
- 29 [34][32] 5-bromo-1-butyl-N-(1-carbamoyl-2,2-dimethyl-propyl)indazole-3-  
30 carboxamide - Other names: ADB-5'Br-BINACA and ADB-5'Br-  
31 BUTINACA.



- 1 [33] Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-  
2 carboxamido)butanoate - Other name: MDMB-4en-PINACA.
- 3 [34] Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-  
4 butanoate - Other names: 4F-MDMB-BUTICA; 4F-MDMB-BICA.
- 5 [35] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-en-1-yl)-1H-  
6 indazole-3-carboxamide - Other name: ADB-4en-PINACA.
- 7 [36] Ethyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-dimethyl-  
8 butanoate - Other names: 5F-EDMB-PICA; 5F-EDMB-2201.
- 9 [37] Methyl 2-(1-(4-fluorobenzyl)-1H-indole-3-carboxamido)-3-methyl  
10 butanoate - Other name: MMB-FUBICA.
- 11 [38] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1H-indazole-3-  
12 carboxamide - Other name: ADB-BUTINACA.
- 13 (4) Indole carboxylic acids. Any compound structurally derived from 1H-indole-  
14 3-carboxylic acid or 1H-2-carboxylic acid substituted in both of the following  
15 ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,  
16 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-  
17 piperidiny)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidiny)methyl,  
18 1-(N-methyl-3-morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo  
19 benzyl group; and, at the hydroxyl group of the carboxylic acid by a phenyl,  
20 benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, or propionaldehyde group  
21 whether or not the compound is further modified to any extent in the  
22 following ways:
- 23 (a) Substitution to the indole ring to any extent; or
- 24 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,  
25 cyclopropyl, propionaldehyde group to any extent; or
- 26 (c) A nitrogen heterocyclic analog of the indole ring; or
- 27 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,  
28 adamantyl, or cyclopropyl ring.
- 29 (e) Examples include:
- 30 [1] 1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid 8-quinolinyl  
31 ester - Other names: BB-22 and QUCHIC.

- 1 [2] naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate -  
2 Other names: FDU-PB-22.
- 3 [3] 1-pentyl-1H-indole-3-carboxylic acid 8-quinolinyl ester - Other  
4 names: PB-22 and QUPIC.
- 5 [4] 1-(5-Fluoropentyl)-1H-indole-3-carboxylic acid 8-quinolinyl ester -  
6 Other names: 5-Fluoro PB-22 and 5F-PB-22.
- 7 [5] quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate - Other  
8 names: FUB-PB-22.
- 9 [6] naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate -  
10 Other names: NM2201 and CBL2201.
- 11 (5) Naphthylmethylindoles. Any compound containing a 1H-indol-3-yl-(1-  
12 naphthyl)methane structure with substitution at the nitrogen atom of the  
13 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,  
14 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-  
15 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or  
16 (tetrahydropyran-4-yl)methyl group whether or not further substituted in the  
17 indole ring to any extent and whether or not substituted in the naphthyl ring  
18 to any extent. Examples include:
- 19 (a) 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane - Other names: JWH-175.  
20 (b) 1-Pentyl-1H-indol-3-yl-(4-methyl-1-naphthyl)methane - Other names:  
21 JWH-184.
- 22 (6) Naphthoylpyrroles. Any compound containing a 3-(1-naphthoyl)pyrrole  
23 structure with substitution at the nitrogen atom of the pyrrole ring by an  
24 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-  
25 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-  
26 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-  
27 yl)methyl group whether or not further substituted in the pyrrole ring to any  
28 extent, whether or not substituted in the naphthyl ring to any extent.  
29 Examples include: (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-  
30 ylmethanone - Other names: JWH-307.

- 1 (7) Naphthylmethylindenes. Any compound containing a naphthylideneindene  
2 structure with substitution at the 3-position of the indene ring by an alkyl,  
3 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-  
4 2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-  
5 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-  
6 yl)methyl group whether or not further substituted in the indene ring to any  
7 extent, whether or not substituted in the naphthyl ring to any extent.  
8 Examples include: E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane  
9 - Other names: JWH-176.
- 10 (8) Cyclohexylphenols. Any compound containing a 2-(3-  
11 hydroxycyclohexyl)phenol structure with substitution at the 5-position of the  
12 phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,  
13 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-  
14 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or  
15 (tetrahydropyran-4-yl)methyl group whether or not substituted in the  
16 cyclohexyl ring to any extent. Examples include:
- 17 (a) 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other  
18 names: CP 47,497.
- 19 (b) 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other  
20 names: Cannabicyclohexanol and CP 47,497 C8 homologue.
- 21 (c) 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-  
22 hydroxypropyl)cyclohexyl]-phenol - Other names: CP 55,940.
- 23 (9) Others specifically named:
- 24 (a) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-  
25 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names: HU-210.
- 26 (b) (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-  
27 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names:  
28 Dexanabinol and HU-211.
- 29 (c) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-  
30 benzoxazin-6-yl]-1-naphthalenylmethanone - Other names:  
31 WIN 55,212-2.

- 1 (d) Naphthalen-1-yl-(4-pentyloxynaphthalen-1-yl)methanone - Other  
2 names: CB-13.
- 3 (e) N-[(Z)-(1-hexyl-2-oxo-indolin-3-ylidene)amino]benzamide - Other  
4 names: BZO-HEXOXIZID and MDA-19.
- 5 (f) N-[(Z)-(2-oxo-1-pentyl-indolin-3-ylidene)amino]benzamide - Other  
6 names: BZO-POXIZID, Pentyl MDA-19, and 5C-MDA-19.
- 7 (g) N-[(Z)-[1-(5-fluoropentyl)-2-oxo-indolin-3-ylidene]amino]benzamide -  
8 Other names: 5F-BZO-POXIZID and 5F-MDA-19.
- 9 (h) N-[(Z)-(2-oxo-1-pent-4-enyl-indolin-3-ylidene)amino]benzamide -  
10 Other names: BZO-4en-POXIZID and 4en-pentyl MDA-19.
- 11 (i) N-[(Z)-[1-(cyclohexylmethyl)-2-oxo-indolin-3-ylidene]amino]benzamide  
12 - Other names: BZO-CHMOXIZID, Cyclohexylmethyl MDA-19 and  
13 CHM-MDA-19.
- 14 (j) N-(1-carbamoyl-2-methyl-propyl)-2-(5-fluoropentyl)-5-(4-  
15 fluorophenyl)pyrazole-3-carboxamide - Other Names: 5F-AB-  
16 PFUPPYCA.
- 17 (k) 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido[4,3-b]indol-1-one - Other  
18 names: CUMYL-PEGACLONE; SGT-151.
- 19 o. Substituted phenethylamines. This includes any compound, unless specifically  
20 excepted, specifically named in this schedule, or listed under a different  
21 schedule, structurally derived from phenylethan-2-amine by substitution on the  
22 phenyl ring in any of the following ways, that is to say, by substitution with a fused  
23 methylenedioxy ring, fused furan ring, or fused tetrahydrofuran ring; by  
24 substitution with two alkoxy groups; by substitution with one alkoxy and either  
25 one fused furan, tetrahydrofuran, or tetrahydropyran ring system; or by  
26 substitution with two fused ring systems from any combination of the furan,  
27 tetrahydrofuran, or tetrahydropyran ring systems.
- 28 (1) Whether or not the compound is further modified in any of the following  
29 ways, that is to say:
- 30 (a) By substitution of phenyl ring by any halo, hydroxyl, alkyl,  
31 trifluoromethyl, alkoxy, or alkylthio groups;

- 1 (b) By substitution at the 2-position by any alkyl groups; or
- 2 (c) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl,
- 3 hydroxybenzyl, methylenedioxybenzyl, or methoxybenzyl groups.
- 4 (2) Examples include:
- 5 (a) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (also known as 2C-C or
- 6 2,5-Dimethoxy-4-chlorophenethylamine).
- 7 (b) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (also known as 2C-D or
- 8 2,5-Dimethoxy-4-methylphenethylamine).
- 9 (c) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (also known as 2C-E or
- 10 2,5-Dimethoxy-4-ethylphenethylamine).
- 11 (d) 2-(2,5-Dimethoxyphenyl)ethanamine (also known as 2C-H or 2,5-
- 12 Dimethoxyphenethylamine).
- 13 (e) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (also known as 2C-I or
- 14 2,5-Dimethoxy-4-iodophenethylamine).
- 15 (f) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (also known as 2C-N or
- 16 2,5-Dimethoxy-4-nitrophenethylamine).
- 17 (g) 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (also known as 2C-
- 18 P or 2,5-Dimethoxy-4-propylphenethylamine).
- 19 (h) 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (also known as 2C-
- 20 T-2 or 2,5-Dimethoxy-4-ethylthiophenethylamine).
- 21 (i) 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (also known as
- 22 2C-T-4 or 2,5-Dimethoxy-4-isopropylthiophenethylamine).
- 23 (j) 2-(4-bromo-2,5-dimethoxyphenyl)ethanamine (also known as 2C-B or
- 24 2,5-Dimethoxy-4-bromophenethylamine).
- 25 (k) 2-(2,5-dimethoxy-4-(methylthio)phenyl)ethanamine (also known as
- 26 2C-T or 4-methylthio-2,5-dimethoxyphenethylamine).
- 27 (l) 1-(2,5-dimethoxy-4-iodophenyl)-propan-2-amine (also known as DOI
- 28 or 2,5-Dimethoxy-4-iodoamphetamine).
- 29 (m) 1-(4-Bromo-2,5-dimethoxyphenyl)-2-aminopropane (also known as
- 30 DOB or 2,5-Dimethoxy-4-bromoamphetamine).

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

- 1 (aa) 5-(2-Aminopropyl)-2,3-dihydrobenzofuran (also known as 5-APDB).
- 2 (bb) 6-(2-Aminopropyl)-2,3,-dihydrobenzofuran (also known as 6-APDB).
- 3 (cc) 2,5-dimethoxy-amphetamine (also known as 2,5-dimethoxy-a-
- 4 methylphenethylamine; 2,5-DMA).
- 5 (dd) 2,5-dimethoxy-4-ethylamphetamine (also known as DOET).
- 6 (ee) 2,5-dimethoxy-4-(n)-propylthiophenethylamine (also known as 2C-T-
- 7 7).
- 8 (ff) 5-methoxy-3,4-methylenedioxy-amphetamine.
- 9 (gg) 4-methyl-2,5-dimethoxy-amphetamine (also known as 4-methyl-2,5-
- 10 dimethoxy-a-methylphenethylamine; DOM and STP).
- 11 (hh) 3,4-methylenedioxy amphetamine (also known as MDA).
- 12 (ii) 3,4-methylenedioxymethamphetamine (also known as MDMA).
- 13 (jj) 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-
- 14 alpha-methyl-3,4(methylenedioxy)phenethylamine, MDE, MDEA).
- 15 (kk) 3,4,5-trimethoxy amphetamine.
- 16 (ll) Mescaline (also known as 3,4,5-trimethoxyphenethylamine).
- 17 p. Substituted tryptamines. This includes any compound, unless specifically
- 18 excepted, specifically named in this schedule, or listed under a different
- 19 schedule, structurally derived from 2-(1H-indol-3-yl)ethanamine (i.e., tryptamine)
- 20 by mono- or di-substitution of the amine nitrogen with alkyl or alkenyl groups or
- 21 by inclusion of the amino nitrogen atom in a cyclic structure whether or not the
- 22 compound is further substituted at the alpha-position with an alkyl group or
- 23 whether or not further substituted on the indole ring to any extent with any alkyl,
- 24 alkoxy, halo, hydroxyl, or acetoxy groups. Examples include:
- 25 (1) 5-methoxy-N,N-diallyltryptamine (also known as 5-MeO-DALT).
- 26 (2) 4-acetoxy-N,N-dimethyltryptamine (also known as 4-AcO-DMT or O-
- 27 Acetylpsilocin).
- 28 (3) 4-hydroxy-N-methyl-N-ethyltryptamine (also known as 4-HO-MET).
- 29 (4) 4-hydroxy-N,N-diisopropyltryptamine (also known as 4-HO-DIPT).
- 30 (5) 5-methoxy-N-methyl-N-isopropyltryptamine (also known as 5-MeO-MiPT).
- 31 (6) 5-methoxy-N,N-dimethyltryptamine (also known as 5-MeO-DMT).

- 1 (7) Bufotenine (also known as 3-(Beta-Dimethyl-aminoethyl)-5-hydroxyindole;  
2 3-(2-dimethylaminoethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-  
3 dimethyltryptamine; mappine).
- 4 (8) 5-methoxy-N,N-diisopropyltryptamine (also known as 5-MeO-DiPT).
- 5 (9) Diethyltryptamine (also known as N,N-Diethyltryptamine; DET).
- 6 (10) Dimethyltryptamine (also known as DMT).
- 7 (11) Psilocyn.
- 8 q. 1-[3-(trifluoromethylphenyl)]piperazine (also known as TFMPP).
- 9 r. 1-[4-(trifluoromethylphenyl)]piperazine.
- 10 s. 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (also known as 5,6-  
11 Methylenedioxy-2-aminoindane or MDAI).
- 12 t. 2-(Ethylamino)-2-(3-methoxyphenyl)cyclohexanone (also known as  
13 Methoxetamine or MXE).
- 14 u. Ethylamine analog of phencyclidine (also known as N-ethyl-1-  
15 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl)  
16 ethylamine, cyclohexamine, PCE).
- 17 v. Pyrrolidine analog of phencyclidine (also known as 1-(1-phenylcyclohexyl)-  
18 pyrrolidine, PCPy, PHP).
- 19 w. Thiophene analog of phencyclidine (also known as (1-[1-(2-thienyl) cyclohexyl]  
20 piperidine; 2-Thienyl analog of phencyclidine; TPCP, TCP).
- 21 x. 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (also known as TCPy).
- 22 y. Salvia divinorum, salvinorin A, or any of the active ingredients of salvia divinorum.
- 23 6. Depressants. Unless specifically excepted or unless listed in another schedule, any  
24 material compound, mixture, or preparation which contains any quantity of the  
25 following substances having a depressant effect on the central nervous system,  
26 whenever the existence of such salts, isomers, and salts of isomers is possible within  
27 the specific chemical designation:
- 28 a. Gamma-hydroxybutyric acid.
- 29 b. Mecloqualone.
- 30 c. Methaqualone.
- 31 d. Clonazolam (also known as Clonitrazolam).



1 e. Etizolam.

2 f. Flualprazolam.

3 g. Flubromazepam.

4 h. Flubromazolam.

5 i. Adinazolam.

6 j. Bromazolam.

7 k. Deschloroetizolam.

8 l. Diclazepam.

9 7. Stimulants. Unless specifically excepted or unless listed in another schedule, any  
10 material, compound, mixture, or preparation which contains any quantity of the  
11 following substances having a stimulant effect on the central nervous system,  
12 including its salts, isomers, and salts of isomers:

13 a. Aminorex (also known as 2-amino-5-phenyl-2-oxazoline, or 4,5-dihydro-5-phenyl-  
14 2-oxazolamine).

15 b. Cathinone.

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

22 (1) By substitution in the ring system to any extent with alkyl, alkylendioxy,  
23 alkoxy, haloalkyl, hydroxyl, or halide substituents, whether or not further  
24 substituted in the ring system by one or more other univalent substituents;

25 (2) By substitution at the 3-position with an acyclic alkyl substituent;

26 (3) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or  
27 methoxybenzyl groups; or

28 (4) By inclusion of the 2-amino nitrogen atom in a cyclic structure.

29 Some trade or other names:

30 (a) 3,4-Methylenedioxy-alpha-pyrrolidinopropiophenone (also known as  
31 MDP PP).

- 1 (b) 3,4-Methylenedioxy-N-ethylcathinone (also known as Ethylone,  
2 MDEC, or bk-MDEA).
- 3 (c) 3,4-Methylenedioxy-N-methylcathinone (also known as Methylone or  
4 bk-MDMA).
- 5 (d) 3,4-Methylenedioxypyrovalerone (also known as MDPV).
- 6 (e) 3,4-Dimethylmethcathinone (also known as 3,4-DMMC).
- 7 (f) 2-(methylamino)-1-phenylpentan-1-one (also known as Pentedrone).
- 8 (g) 2-Fluoromethcathinone (also known as 2-FMC).
- 9 (h) 3-Fluoromethcathinone (also known as 3-FMC).
- 10 (i) 4-Methylethcathinone (also known as 4-MEC and 4-methyl-N-  
11 ethylcathinone).
- 12 (j) 4-Fluoromethcathinone (also known as Flephedrone and 4-FMC).
- 13 (k) 4-Methoxy-alpha-pyrrolidinopropiophenone (also known as MOPPP).
- 14 (l) 4-Methoxymethcathinone (also known as Methedrone; bk-PMMA).
- 15 (m) 4'-Methyl-alpha-pyrrolidinobutiophenone (also known as MPBP).
- 16 (n) Alpha-methylamino-butyrophenone (also known as Buphedrone or  
17 MABP).
- 18 (o) Alpha-pyrrolidinobutiophenone (also known as alpha-PBP).
- 19 (p) Alpha-pyrrolidinopropiophenone (also known as alpha-PPP).
- 20 (q) Alpha-pyrrolidinopentiophenone (also known as Alpha-  
21 pyrrolidinovalerophenone or alpha-PVP).
- 22 (r) Beta-keto-N-methylbenzodioxolylbutanamine (also known as Butylone  
23 or bk-MBDB).
- 24 (s) Ethcathinone (also known as N-Ethylcathinone).
- 25 (t) 4-Methylmethcathinone (also known as Mephedrone or 4-MMC).
- 26 (u) Methcathinone.
- 27 (v) N,N-dimethylcathinone (also known as metamfepramone).
- 28 (w) Naphthylpyrovalerone (naphyrone).
- 29 (x) B-Keto-Methylbenzodioxolylpentanamine (also known as Pentylone).
- 30 (y) 4-Methyl-alpha-pyrrolidinopropiophenone (also known as 4-MePPP  
31 and MPPP).

- 1 (z) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one (also known as  
2 Ephylone and N-Ethylpentylone).
- 3 (aa) N-ethylhexedrone (also known as alpha - ethylaminohexanophenone  
4 and 2-(ethylamino)-1-phenylhexan-1-one)).
- 5 (bb) Alpha-pyrrolidinohexanophenone (also known as alpha-PHP, alpha-  
6 pyrrolidinohexiophenone, and 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-  
7 one)).
- 8 (cc) 4-methyl-alpha-ethylaminopentiophenone (also known as 4-MEAP  
9 and 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one)).
- 10 (dd) 4'-methyl-alpha-pyrrolidinohexiophenone (also known as MPHP, 4'-  
11 methyl-alpha-pyrrolidinohexanophenone and 1-(4-methylphenyl)-2-  
12 (pyrrolidin-1-yl)hexan-1-one)).
- 13 (ee) Alpha-pyrrolidinoheptaphenone (also known as PV8 and 1-phenyl-2-  
14 (pyrrolidin-1-yl)heptan-1-one)).
- 15 (ff) 4-chloro-alpha-pyrrolidinovalerophenone (also known as 4-chloro-  
16 alpha-PVP, 4'-chloro-alpha-pyrrolidinopentiophenone, and 1-(4-  
17 chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one)).
- 18 (gg) 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one (other name: alpha-  
19 PiHP).
- 20 (hh) 2-(methylamino)-1-(3-methylphenyl)propan-1-one (other names: 3-  
21 MMC; 3-methylmethcathinone).
- 22 (ii) Eutylone (also known as 1-(1,3-benzodioxol-5-yl)-2-  
23 (ethylamino)butan-1-one).
- 24 d. Fenethylline.
- 25 e. Fluoroamphetamine.
- 26 f. Fluoromethamphetamine.
- 27 g. ( $\pm$ )cis-4-methylaminorex (also known as ( $\pm$ )cis-4,5-dihydro-4-methyl-5-phenyl-2-  
28 oxazolamine).
- 29 h. N-Benzylpiperazine (also known as BZP, 1-benzylpiperazine).
- 30 i. N-ethylamphetamine.

- 1           j.    N, N-dimethylamphetamine (also known as N,N-alpha-trimethyl-  
2                    benzeneethanamine; N,N-alpha-trimethylphenethylamine).  
3           k.    1-(4-methoxyphenyl)-N-methylpropan-2-amine (also known as  
4                    paramethoxymethamphetamine and PMMA).  
5           l.    4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-  
6                    oxazolamine; 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine).  
7           m.    Amineptine (Also known as 7- [(10,11-dihydro-5Hdibenzo[a,d]cyclohepten-5-  
8                    yl)amino]heptanoic acid).  
9           n.    Mesocarb (Also known as N-phenyl-N' -(3-(1- phenylpropan-2-yl)-1,2,3-  
10                   oxadiazol-3-ium-5-yl)carbamidate).  
11          o.    Methiopropamine (Also known as N-methyl-1-(thiophen-2-yl)propan-2-amine).  
12          p.    Ethylphenidate (ethyl 2-phenyl-2-(piperidin-2-yl)acetate).

13           **SECTION 2. AMENDMENT.** Section 19-03.1-09 of the North Dakota Century Code is  
14 amended and reenacted as follows:

15           **19-03.1-09. Schedule III.**

- 16           1.    The controlled substances listed in this section are included in schedule III.  
17           2.    Schedule III consists of the drugs and other substances, by whatever official name,  
18                   common or usual name, chemical name, or brand name designated, listed in this  
19                   section.  
20           3.    Stimulants. Unless specifically excepted or unless listed in another schedule, any  
21                   material, compound, mixture, or preparation which contains any quantity of the  
22                   following substances having a stimulant effect on the central nervous system,  
23                   including its salts, isomers (whether optical, position, or geometric), and salts of such  
24                   isomers whenever the existence of such salts, isomers, and salts of isomers is  
25                   possible within the specific chemical designation:  
26                   a.    Those compounds, mixtures, or preparations in dosage unit form containing any  
27                   stimulant substances listed in schedule II and any other drug of the quantitative  
28                   composition shown in that schedule for those drugs or which is the same except  
29                   that it contains a lesser quantity of controlled substances.  
30                   b.    Benzphetamine.  
31                   c.    Chlorphentermine.

- 1           d. Clortermine.
- 2           e. Phendimetrazine.
- 3        4. Depressants. Unless specifically excepted or unless listed in another schedule, any
- 4           material, compound, mixture, or preparation that contains any quantity of the following
- 5           substances having a depressant effect on the central nervous system:
- 6           a. Any compound, mixture, or preparation containing:
- 7                (1) Amobarbital;
- 8                (2) Secobarbital;
- 9                (3) Pentobarbital;
- 10           or any salt thereof and one or more other active medicinal ingredients which are
- 11           not listed in any schedule.
- 12           b. Any suppository dosage form containing:
- 13                (1) Amobarbital;
- 14                (2) Secobarbital;
- 15                (3) Pentobarbital;
- 16           or any salt of any of these drugs and approved by the food and drug
- 17           administration for marketing only as a suppository.
- 18           c. Any substance that contains any quantity of a derivative of barbituric acid, or any
- 19           salt of a derivative of barbituric acid, except those substances which are
- 20           specifically listed in other schedules thereof.
- 21           d. Chlorhexadol.
- 22           e. Embutramide.
- 23           f. Gamma-hydroxybutyric acid in a United States food and drug administration-
- 24           approved drug product.
- 25           g. Ketamine.
- 26           h. Lysergic acid.
- 27           i. Lysergic acid amide.
- 28           j. Methyprylon.
- 29           k. Perampanel.
- 30           l. Sativex or its successor name as determined by the federal food and drug
- 31           administration.

- 1 m. Sulfondiethylmethane.
- 2 n. Sulfonethylmethane.
- 3 o. Sulfonmethane.
- 4 p. Tiletamine and zolazepam or any salt thereof. Some trade or other names for a  
5 tiletamine-zolazepam combination product: Telazol. Some trade or other names  
6 for tiletamine: 2-(ethylamino)-2-(2-thienyl)-cyclohexanone. Some trade or other  
7 names for zolazepam: 4-2(2-fluorophenyl)-6, 8-dihydro-1,3,8-trimethylpyrazolo-  
8 [3,4-e][1,4]-diazepin-7(1H)-one, flupyrzapon.
- 9 5. Nalorphine.
- 10 6. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any  
11 material, compound, mixture, or preparation that contains any of the following narcotic  
12 drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited  
13 quantities as set forth below:
  - 14 a. (1) Not more than 1.80 grams of codeine per 100 milliliters or not more than  
15 90 milligrams per dosage unit, with an equal or greater quantity of an  
16 isoquinoline alkaloid of opium.
  - 17 (2) Not more than 1.80 grams of codeine per 100 milliliters or not more than  
18 90 milligrams per dosage unit, with one or more active, non-narcotic  
19 ingredients in recognized therapeutic amounts.
  - 20 (3) Not more than 1.80 grams of dihydrocodeine per 100 milliliters or not more  
21 than 90 milligrams per dosage unit, with one or more active, non-narcotic  
22 ingredients in recognized therapeutic amounts.
  - 23 (4) Not more than 300 milligrams of ethylmorphine per 100 milliliters or not  
24 more than 15 milligrams per dosage unit, with one or more active,  
25 non-narcotic ingredients in recognized therapeutic amounts.
  - 26 (5) Not more than 500 milligrams of opium per 100 milliliters or per 100 grams,  
27 or not more than 25 milligrams per dosage unit, with one or more active,  
28 non-narcotic ingredients in recognized therapeutic amounts.
  - 29 (6) Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams  
30 with one or more active, non-narcotic ingredients in recognized therapeutic  
31 amounts.

- 1           b. Buprenorphine.
- 2           7. Anabolic steroids. Unless specifically excepted or unless listed in another schedule,  
3           any material, compound, mixture, or preparation that contains any quantity of the  
4           following ~~anabolic steroids~~ substances, including its salts, esters, and ethers:
- 5           a. 3beta,17-dihydroxy-5a-androstane;
- 6           b. 3alpha,17beta-dihydroxy-5a-androstane;
- 7           c. 5alpha-androstan-3,17-dione;
- 8           d. 5alpha-androstan-3,6,17-trione;
- 9           e. 1-androstenediol (3beta,17beta-dihydroxy-5alpha-androst-1-ene);
- 10          e-f. 1-androstenediol (3alpha,17beta-dihydroxy-5alpha-androst-1-ene);
- 11          f-g. 4-androstenediol (3beta,17beta-dihydroxyandrost-4-ene);
- 12          g-h. 5-androstenediol (3beta,17beta-dihydroxy-androst-5-ene);
- 13          h-i. 1-androstenedione ([5alpha]-androst-1-en-3,17-dione);
- 14          i-j. 4-androstenedione (androst-4-en-3,17-dione);
- 15          j-k. 5-androstenedione (androst-5-en-3,17-dione);
- 16          k-l. Bolasterone (7alpha,17alpha-dimethyl-17beta-hydroxyandrost-4-en-3-one);
- 17          l-m. Boldenone (17beta-hydroxyandrost-1,4,-diene-3-one);
- 18          m-n. Boldione (androsta-1,4-diene-3,17-dione);
- 19          o. 6-bromo-androsta-1,4-diene-3,17-dione;
- 20          p. 6-bromo-androstan-3,17-dione;
- 21          n-q. Calusterone (7beta,17alpha-dimethyl-17beta-hydroxyandrost-4-en-3-one);
- 22          r. 4-chloro-17alpha-methyl-androsta-1,4-diene-3,17beta-diol;
- 23          s. 4-chloro-17alpha-methyl-androst-4-ene-3beta,17beta-diol;
- 24          t. 4-chloro-17alpha-methyl-17beta-hydroxy-androst-4-en-3-one;
- 25          u. 4-chloro-17alpha-methyl-17beta-hydroxy-androst-4-ene-3,11-dione;
- 26          e-v. Clostebol (4-chloro-17beta-hydroxyandrost-4-en-3-one);
- 27          p-w. Dehydrochloromethyltestosterone (4-chloro-17beta-hydroxy-17alpha-methyl-  
28          androst-1,4-dien-3-one);
- 29          q-x. Delta-1-dihydrotestosterone (also known as '1-testosterone') (17beta-hydroxy-  
30          5alpha-androst-1-en-3-one);

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- 1           r.y.   Desoxymethyltestosterone (17a-methyl-5a-androst-2-en-17beta-ol) (also known  
2           as madol);
- 3           s.z.   4-dihydrotestosterone (17beta-hydroxy-androstan-3-one);
- 4           aa.   3beta,17beta-dihydroxy-5alpha-androstane;
- 5           bb.   3alpha,17beta-dihydroxy-5alpha-androstane;
- 6           cc.   2alpha,17alpha-dimethyl-17beta-hydroxy-5beta-androstan-3-one;
- 7           t.dd.   Drostanolone (17beta-hydroxy-2alpha-methyl-5alpha-androstan-3-one);
- 8           ee.   2alpha,3alpha-epithio-17alpha-methyl-5alpha-androstan-17beta-ol;
- 9           ff.   estra-4,9,11-triene-3,17-dione;
- 10          gg.   13beta-ethyl-17beta-hydroxygon-4-en-3-one;
- 11          u.hh.   Ethylestrenol (17alpha-ethyl-17beta-hydroxyestr-4-ene);
- 12          v.ii.   Fluoxymesterone (9-fluoro-17alpha-methyl-11beta, 17beta-dihydroxyandrost-4-  
13          en-3-one);
- 14          w.jj.   Formebolone (2-formyl-17alpha-methyl-11alpha, 17beta-dihydroxyandrost-1,4-  
15          dien-3-one);
- 16          x.kk.   Furazabol (17alpha-methyl-17beta-hydroxyandrostan[2,3-c]-furazan);
- 17          ll.   [3,2-c]furazan-5alpha-androstan-17beta-ol;
- 18          mm.   18a-homo-3-hydroxy-estra-2,5(10)-dien-17-one;
- 19          nn.   4-hydroxy-androst-4-ene-3,17-dione;
- 20          oo.   17beta-hydroxy-androstano[2,3-d]isoxazole;
- 21          pp.   17beta-hydroxy-androstano[3,2-c]isoxazole;
- 22          qq.   3beta-hydroxy-estra-4,9,11-trien-17-one;
- 23          y.rr.   13beta-ethyl-17alpha-hydroxygon-4-en-3-one;
- 24          z.ss.   4-hydroxytestosterone (4,17beta-dihydroxy-androst-4-en-3-one);
- 25          aa.tt.   4-hydroxy-19-nortestosterone (4,17beta-dihydroxy-estr-4-en-3-one);
- 26          bb.uu.   Mestanolone (17alpha-methyl-17beta-hydroxy-5alpha-androstan-3-one);
- 27          ee.vv.   Mesterolone (1alpha-methyl-17beta-hydroxy-[5alpha]-androstan-3-one);
- 28          dd.ww.   Methandienone (17alpha-methyl-17beta-dihydroxyandrost-1,4-dien-3-one);
- 29          ee.xx.   Methandriol (17alpha-methyl-3beta,17beta-dihydroxyandrost-5-ene);
- 30          ff.yy.   Methasterone (2[alpha],17[alpha]-dimethyl-5[alpha]-androstan-17[beta]-ol-3-one);
- 31          gg.zz.   Methenolone (1-methyl-17beta-hydroxy-5alpha-androst-1-en-3-one);



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- 1        aaa.    17alpha-methyl-androsta-1,4-diene-3,17beta-diol:
- 2        bbb.    17alpha-methyl-5alpha-androstan-17beta-ol:
- 3        ccc.    17alpha-methyl-androstan-3-hydroxyimine-17beta-ol:
- 4        ddd.    6alpha-methyl-androst-4-ene-3,17-dione;
- 5        eee.    17alpha-methyl-androst-2-ene-3,17beta-diol:
- 6        hh.fff. 17alpha-methyl-3beta,17beta-dihydroxy-5a-androstane;
- 7        ii.ggg. 17alpha-methyl-3alpha,17beta-dihydroxy-5a-androstane;
- 8        jj.hhh. 17alpha-methyl-3beta,17beta-dihydroxyandrost-4-ene;
- 9        kk.iii. 17alpha-methyl-4-hydroxynandrolone (17alpha-methyl-4-hydroxy-17beta-
- 10       hydroxyestr-4-en-3-one);
- 11       ll.jjj. Methyldienolone (17alpha-methyl-17beta-hydroxyestra-4,9(10)-dien-3-one);
- 12       mm.kkk.        Methyltrienolone (17alpha-methyl-17beta-hydroxyestra-4,9(11)-trien-3-
- 13       one);
- 14       nn.lll. Methyltestosterone (17alpha-methyl-17beta-hydroxyandrost-4-en-3-one);
- 15       oo.mmm.        Mibolerone (7alpha,17alpha-dimethyl-17beta-hydroxyestr-4-en-3-one);
- 16       pp.nnn. 17alpha-methyl-delta1-dihydrotestosterone (17beta-hydroxy-17alpha-methyl-
- 17       5alpha-androst-1-en-3-one) (also known as '17-alpha-methyl-1-testosterone');
- 18       qq.ooo. Nandrolone (17beta-hydroxyestr-4-en-3-one);
- 19       rr.ppp. 19-nor-4-androstenediol (3beta,17beta-dihydroxyestr-4-ene);
- 20       ss.ggg. 19-nor-4-androstenediol (3alpha,17beta-dihydroxyestr-4-ene);
- 21       tt.rrr. 19-nor-5-androstenediol (3beta,17beta-dihydroxyestr-5-ene);
- 22       uu.sss. 19-nor-5-androstenediol (3alpha,17beta-dihydroxyestr-5-ene);
- 23       vv.ttt. 19-nor-4-androstenedione (estr-4-en-3,17-dione);
- 24       ww.uuu.        19-nor-4,9(10)-androstadienedione (estra-4,9(10)-diene-3,17-dione);
- 25       xx.vvv. 19-nor-5-androstenedione (estr-5-en-3,17-dione);
- 26       yy.www. Norbolethone (13beta,17alpha-diethyl-17beta-hydroxygon-4-en-3-one);
- 27       zz.xxx. Norclostebol (4-chloro-17beta-hydroxyestr-4-en-3-one);
- 28       aaa.yyy.        Norethandrolone (17alpha-ethyl-17beta-hydroxyestr-4-en-3-one);
- 29       bbb.zzz.        Normethandrolone (17alpha-methyl-17beta-hydroxyestr-4-en-3-one);
- 30       eee.aaaa.       Oxandrolone (17alpha-methyl-17beta-hydroxy-2-oxa-[5alpha]-androstan-
- 31       3-one);

- 1        ddd.bbbb.     Oxymesterone (17alpha-methyl-4-17beta-dihydroxyandrost-4-en-3-one);
- 2        eee.cccc.     Oxymetholone (17alpha-methyl-2-hydroxymethylene-17beta-hydroxy
- 3                    [5alpha]-androstan-3-one);
- 4        dddd.    [3,2-c]pyrazole-androst-4-en-17beta-ol;
- 5        fff.eeee. Stanazolol (17alpha-methyl-17beta-hydroxy[5alpha]-androst-2-eno[3,2-c]-
- 6                    pyrazole);
- 7        ggg.ffff. Stenbolone (17beta-hydroxy-2-methyl-[5alpha]-androst-1-en-3-one);
- 8        hhh.gggg.     Prostanazol (17[beta]- hydroxy-5[alpha]-androstano[3,2-c]pyrazole);
- 9        iii.hhhh. Testolactone (13-hydroxy-3-oxo-13,17-secoandrosta-1,4-dien-17-oic acid
- 10                   lactone);
- 11        jjj.iiii.     Testosterone (17beta-hydroxyandrost-4-en-3-one);
- 12        kkk.jjjj. Tetrahydrogestrinone (13beta,17alpha-diethyl-17beta-hydroxygon-4,9,11-trien-3-
- 13                    one); or
- 14        ###kkkk. Trenbolone (17beta-hydroxyestr-4,9,11-trien-3-one);
- 15                    ~~or any salt, ester, or isomer of a drug or substance described or listed in this~~
- 16                    ~~subsection, if that salt, ester, or isomer promotes muscle growth.~~

17        The term does not include an anabolic steroid that is expressly intended for

18        administration through implants to cattle or other nonhuman species and which has

19        been approved by the secretary of health and human services for administration

20        unless any person prescribes, dispenses, possesses, delivers, or distributes for

21        human use.

22        8. Hallucinogenic substances.

- 23        a. Dronabinol (synthetic) [(-)-delta-9-(trans)-tetrahydrocannabinol] in sesame oil and
- 24                    encapsulated in a soft gelatin capsule in a United States food and drug
- 25                    administration-approved drug product.
- 26        b. Any product in hard or soft gelatin capsule form containing natural dronabinol
- 27                    (derived from the cannabis plant) or synthetic dronabinol (produced from
- 28                    synthetic materials) in sesame oil, for which an abbreviated new drug application
- 29                    has been approved by the food and drug administration under section 505(j) of
- 30                    the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355(j)] which references as
- 31                    its listed drug the drug product referred to in subdivision a.

- 1           9.    The board may except by rule any compound, mixture, or preparation containing any  
2           stimulant or depressant substance listed in subsections 3 and 4 from the application of  
3           all or any part of this chapter if the compound, mixture, or preparation contains one or  
4           more active medicinal ingredients not having a stimulant or depressant effect on the  
5           central nervous system, and if the admixtures are included therein in combinations,  
6           quantity, proportion, or concentration that vitiate the potential for abuse of the  
7           substances which have a stimulant or depressant effect on the central nervous  
8           system.

9           **SECTION 3. AMENDMENT.** Section 19-03.1-11 of the North Dakota Century Code is  
10          amended and reenacted as follows:

11          **19-03.1-11. Schedule IV.**

- 12          1.    The controlled substances listed in this section are included in schedule IV.  
13          2.    Schedule IV consists of the drugs and other substances, by whatever official name,  
14          common or usual name, chemical name, or brand name designated, listed in this  
15          section.  
16          3.    Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any  
17          material, compound, mixture, or preparation containing any of the following narcotic  
18          drugs or their salts calculated as the free anhydrous base or alkaloid, in limited  
19          quantities as set forth below:  
20                  a.    Not more than 1 milligram of difenoxin and not less than 25 micrograms of  
21                  atropine sulfate per dosage unit.  
22                  b.    Dextropropoxyphene (also known as alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-  
23                  methyl-2-propionoxybutane).  
24                  c.    2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its salts, optical  
25                  and geometric isomers and salts of these isomers including Tramadol.  
26          4.    Depressants. Unless specifically excepted or unless listed in another schedule, any  
27          material, compound, mixture, or preparation containing any quantity of the following  
28          substances, including their salts, isomers, and salts of isomers whenever the  
29          existence of those salts, isomers, and salts of isomers is possible within the specific  
30          chemical designation:  
31                  a.    Alprazolam.

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- 1 b. Alfaxalone.
- 2 c. Barbital.
- 3 d. Brexanolone.
- 4 e. Bromazepam.
- 5 f. Camazepam.
- 6 g. Carisoprodol.
- 7 h. Chloral betaine.
- 8 i. Chloral hydrate.
- 9 j. Chlordiazepoxide.
- 10 k. Clobazam.
- 11 l. Clonazepam.
- 12 m. Clorazepate.
- 13 n. Clotiazepam.
- 14 o. Cloxazolam.
- 15 p. Daridorexant.
- 16 q. Delorazepam.
- 17 r. Diazepam.
- 18 s. Dichloralphenazone.
- 19 t. Estazolam.
- 20 u. Ethchlorvynol.
- 21 v. Ethinamate.
- 22 w. Ethyl loflazepate.
- 23 x. Fludiazepam.
- 24 y. Flunitrazepam.
- 25 z. Flurazepam.
- 26 aa. Fospropofol.
- 27 bb. Halazepam.
- 28 cc. Haloxazolam.
- 29 dd. Indiplon.
- 30 ee. Ketazolam.
- 31 ff. Lemborexant.

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1	gg.	Loprazolam.
2	hh.	Lorazepam.
3	ii.	Lorcaserin.
4	jj.	Lormetazepam.
5	kk.	Mebutamate.
6	ll.	Medazepam.
7	mm.	Meprobamate.
8	nn.	Methohexital.
9	oo.	Methylphenobarbital (also known as mephobarbital).
10	pp.	Midazolam.
11	qq.	Nimetazepam.
12	rr.	Nitrazepam.
13	ss.	Nordiazepam.
14	tt.	Oxazepam.
15	uu.	Oxazolam.
16	vv.	Paraldehyde.
17	ww.	Petrichloral.
18	xx.	Phenobarbital.
19	yy.	Pinazepam.
20	zz.	Propofol.
21	aaa.	Prazepam.
22	bbb.	Quazepam.
23	ccc.	Remimazolam.
24	ddd.	Suvorexant.
25	eee.	Temazepam.
26	fff.	Tetrazepam.
27	ggg.	Triazolam.
28	hhh.	Zaleplon.
29	iii.	Zolpidem.
30	jjj.	Zopiclone.
31	<u>kkk.</u>	<u>Zuranolone.</u>

- 1           5. Stimulants. Unless specifically excepted or unless listed in another schedule, any  
2           material, compound, mixture, or preparation which contains any quantity of the  
3           following substances having a stimulant effect on the central nervous system,  
4           including its salts, isomers, and salts of isomers:
- 5           a. Cathine.  
6           b. Diethylpropion.  
7           c. Fencamfamin.  
8           d. Fenproporex.  
9           e. Mazindol.  
10          f. Mefenorex.  
11          g. Modafinil.  
12          h. Pemoline (including organometallic complexes and chelates thereof).  
13          i. Phentermine.  
14          j. Pipradrol.  
15          k. Serdexmethylphenidate.  
16          l. Sibutramine.  
17          m. Solriamfetol.  
18          n. SPA ((-)-1-dimethylamino-1, 2-diphenylethane).
- 19          6. Other substances. Unless specifically excepted or unless listed in another schedule,  
20          any material, compound, mixture, or preparation which contains any quantity of:
- 21          a. Pentazocine, including its salts.  
22          b. Butorphanol, including its optical isomers.  
23          c. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-  
24          oxopropyl]][(1S)-1-(4-phenyl-1*H*-imidazol-2-yl)ethyl]amino]methyl]-2-  
25          methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and  
26          salts of isomers.
- 27          7. Hallucinogenic substances. Pharmaceutical composition of crystalline polymorph  
28          psilocybin, known as COMP360 or any such trade name approved for COMP360 by  
29          the United States food and drug administration.
- 30          8. The board may except by rule any compound, mixture, or preparation containing any  
31          depressant substance listed in subsection 2 from the application of all or any part of

1           this chapter if the compound, mixture, or preparation contains one or more active  
2           medicinal ingredients not having a depressant effect on the central nervous system,  
3           and if the admixtures are included therein in combinations, quantity, proportion, or  
4           concentration that vitiate the potential for abuse of the substances which have a  
5           depressant effect on the central nervous system.

6           **SECTION 4. EMERGENCY.** This Act is declared to be an emergency measure.