

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.
23 h. Betameprodine.
24 i. Betamethadol.

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

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

- 1 iii. 2-(4-ethoxybenzyl)-5-nitro-1-(2-(piperidin-1-yl)ethyl)-1H-benzimidazole (Other
2 names: N-piperidinyl etonitazene; etonitazepipne).
- 3 iii. 2-Methyl AP-237 (1-(2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl)butan-1-
4 one).
- 5 kkk. Fentanyl derivatives. Unless specifically excepted or unless listed in another
6 schedule or are not FDA approved drugs, and are derived from N-(1-(2-
7 Phenylethyl)-4-piperidinyl)-N-phenylpropanamide (Fentanyl) by any substitution
8 on or replacement of the phenethyl group, any substitution on the piperidine ring,
9 any substitution on or replacement of the propanamide group, any substitution on
10 the anilido phenyl group, or any combination of the above. Examples include:
11 (1) N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide (also known
12 as Acetyl-alpha-methylfentanyl).
- 13 (2) N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-
14 2-phenylethyl)-4-(N-propanilido)piperidine (also known as Alpha-
15 methylfentanyl).
- 16 (3) N-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (also
17 known as Alpha-methylthiofentanyl).
- 18 (4) N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide (also
19 known as Beta-hydroxyfentanyl).
- 20 (5) N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide
21 (also known as Beta-hydroxy-3-methylfentanyl).
- 22 (6) N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (also
23 known as 3-Methylfentanyl).
- 24 (7) N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (also
25 known as 3-Methylthiofentanyl).
- 26 (8) N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl]propanamide (also
27 known as Para-fluorofentanyl).
- 28 (9) N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]propanamide (also known as
29 Thiofentanyl).
- 30 (10) N-(1-phenylethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (also known
31 as Furanyl Fentanyl).

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

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

- 1 (41) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-carboxamide (para-
2 fluoro furanyl fentanyl).
- 3 (42) 2',5'-dimethoxyfentanyl(N-(1-(2,5-dimethoxyphenethyl)piperidine-4-yl)-N-
4 phenylpropionamide).
- 5 (43) 3-furanyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3-
6 carboxamide).
- 7 (44) alpha'-methyl butyryl fentanyl(2-methyl-N-(1-phenethylpiperidin-4-yl)-N-
8 phenylbutanamide).
- 9 (45) isovaleryl fentanyl(3-methyl-N-(1-phenethylpiperidin-4-yl)-N-
10 phenylbutanamide).
- 11 (46) meta-fluorofentanyl(N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-
12 yl)propionamide).
- 13 (47) meta-fluoroisobutyryl fentanyl(N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-
14 yl)isobutyramide).
- 15 (48) ortho-fluorofuranyl fentanyl(N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-
16 yl)furan-2-carboxamide).
- 17 (49) para-methoxyfuranyl fentanyl(N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-
18 4-yl)furan-2-carboxamide).
- 19 (50) para-methylcyclopropyl fentanyl(N-(4-methylphenyl)-N-(1-phenylpiperidin-4-
20 yl)cyclopropanecarboxamide).
- 21 4. Opium derivatives. Unless specifically excepted or unless listed in another schedule,
22 any of the following opium derivatives, its salts, isomers, and salts of isomers
23 whenever the existence of such salts, isomers, and salts of isomers is possible within
24 the specific chemical designation:
- 25 a. Acetorphine.
- 26 b. Acetyldihydrocodeine.
- 27 c. Benzylmorphine.
- 28 d. Codeine methylbromide.
- 29 e. Codeine-N-Oxide.
- 30 f. Cyrenorphine.
- 31 g. Desomorphine.

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- 1 h. Dihydromorphine.
- 2 i. Drotebanol.
- 3 j. Etorphine (except hydrochloride salt).
- 4 k. Heroin.
- 5 l. Hydromorphinol.
- 6 m. Methyldesorphine.
- 7 n. Methyldihydromorphine.
- 8 o. Morphine methylbromide.
- 9 p. Morphine methylsulfonate.
- 10 q. Morphine-N-Oxide.
- 11 r. Myrophine.
- 12 s. Nicocodeine.
- 13 t. Nicomorphine.
- 14 u. Normorphine.
- 15 v. Pholcodine.
- 16 w. Thebacon.
- 17 5. Hallucinogenic substances. Unless specifically excepted or unless listed in another
18 schedule, any material, compound, mixture, or preparation containing any quantity of
19 the following hallucinogenic substances, including their salts, isomers, and salts of
20 isomers whenever the existence of those salts, isomers, and salts of isomers is
21 possible within the specific chemical designation (for purposes of this subsection only,
22 the term "isomer" includes the optical, position, and geometric isomers):
 - 23 a. Alpha-ethyltryptamine, its optical isomers, salts, and salts of isomers (also known
24 as etryptamine; α -ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole).
 - 25 b. Alpha-methyltryptamine.
 - 26 c. 4-methoxyamphetamine (also known as 4-methoxy- α -methylphenethylamine;
27 paramethoxyamphetamine; PMA).
 - 28 d. N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy- α -
29 methyl-3,4(methylenedioxy)phenylamine, and N-hydroxy MDA).
 - 30 e. Ibogaine (also known as 7-Ethyl-6, 6B, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-
31 6, 9-methano-5 H-pyrido [1', 2':1,2] azepino (5,4-b) indole; Tabernanthe iboga).

- 1 f. Lysergic acid diethylamide.
- 2 g. Marijuana.
- 3 h. Parahexyl (also known as 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro- 6,6,9-trimethyl-
4 6H-dibenzol[b,d]pyran; Synhexyl).
- 5 i. Peyote (all parts of the plant presently classified botanically as *Lophophora*
6 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from
7 any part of such plant, and every compound, manufacture, salts, derivative,
8 mixture, or preparation of such plant, its seeds, or its extracts).
- 9 j. N-ethyl-3-piperidyl benzilate.
- 10 k. N-methyl-3-piperidyl benzilate.
- 11 l. Psilocybin.
- 12 m. (1) Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained
13 in a plant of the genus *Cannabis* (cannabis plant), as well as synthetic
14 equivalents of the substances contained in the cannabis plant, or in the
15 resinous extractives of such plant, including synthetic substances,
16 derivatives, and their isomers with similar chemical structure and
17 pharmacological activity to those substances contained in the plant; such as
18 the following:
- 19 (a) Delta-1 cis or trans tetrahydrocannabinol, and their optical isomers.
20 Other names: Delta-9-tetrahydrocannabinol.
- 21 (b) Delta-6 cis or trans tetrahydrocannabinol, and their optical isomers.
22 Other names: Delta-8-tetrahydrocannabinol.
- 23 (c) Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers.
24 (Since nomenclature of these substances is not internationally standardized,
25 compounds of these structures, regardless of numerical designation of atomic
26 positions covered.)
- 27 (2) Tetrahydrocannabinols do not include:
- 28 (a) The allowable amount of total tetrahydrocannabinol found in hemp or
29 an allowed hemp commodity or product as defined in chapter
30 4.1-18.1; or

- 1 (b) A prescription drug approved by the United States food and drug
2 administration under section 505 of the Federal Food, Drug, and
3 Cosmetic Act [21 U.S.C. 355].
- 4 n. Cannabinoids, synthetic. It includes the chemicals and chemical groups listed
5 below, including their homologues, salts, isomers, and salts of isomers. The term
6 "isomer" includes the optical, position, and geometric isomers.
- 7 (1) Indole acetamides. Any compound structurally derived from 1H-indole3-
8 acetamide or 1H-2-acetamide substituted in both of the following ways: at
9 the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,
10 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2piperidinyl)methyl, 2-
11 (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 hydrogen of the acetamide by a phenyl, benzyl, cumyl, naphthyl,
14 adamantyl, cyclopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group
15 whether or not the compound is further modified to any extent in the
16 following ways:
- 17 (a) Substitution to the indole ring to any extent; or
18 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
19 cyclopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group to any
20 extent; or
21 (c) A nitrogen heterocyclic analog of the indole ring; or
22 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,
23 adamantyl, or cyclopropyl ring.
24 (e) Examples include:
- 25 [1] N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide - Other names: CH-
26 PIATA, Cyclohexyl-PIATA, CHX-PIATA, CH-PIACA, and CHX-
27 PIACA.
28 [2] N-cyclohexyl-2-[1-[(4-fluorophenyl)methyl]indol-3-yl]acetamide -
29 Other names: CH-FUBIATA and CH-FUBIACA.

- 1 [3] 2-[[2-[1-[(4-fluorophenyl)methyl]indol-3-yl]acetyl]amino]-3,3-
2 dimethyl-butanamide - Other names: ADB-FUBIATA, FUB-
3 ACADB, and AD-18.
- 4 (2) 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 (3) 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] 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-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-
18 carboxamide - Other names: FUB-AKB48, FUB-APINACA, and
19 AKB48 N-(4-FLUOROBENZYL).
- 20 [13] 1-(5-fluoropentyl)-N-(quinolin-8-yl)-1H-indazole-3-carboxamide -
21 Other names: 5-fluoro-THJ.
- 22 [14] methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-
23 methylbutanoate - Other names: 5-fluoro AMB and 5F-AMB.
- 24 [15] methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-
25 methylbutanoate - Other names: FUB-AMB, MMB-FUBINACA,
26 and AMB-FUBINACA.
- 27 [16] N-[1-(aminocarbonyl)-2,2-dimethylpropyl]-1-(cyclohexylmethyl)-1
28 H-indazole-3-carboxamide - Other names: MAB-CHMINACA and
29 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 [24] ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-
21 dimethylbutanoate - Other names: 5F-EDMB-PINACA.
- 22 [25] methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-
23 dimethylbutanoate - Other names: 5F-MDMB-PICA and 5F-
24 MDMB-2201.
- 25 [26] 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-
26 carboxamide - Other names: 5F-CUMYL-PINACA, SGT-25.
- 27 [27] (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)
28 methanone - Other names: FUB-144.
- 29 [28] methyl 2-(1-(4-fluorobutyl)-1H-indazole-3-carboxamido)-3,3-
30 dimethylbutanoate (4F-MDMB-BINACA, 4F-MDMB-BUTINACA).

- 1 [29] Methyl 3,3-dimethyl-2-[(1-pent-4-enyl)indazole-3-
2 carbonyl]amino]butanoate - Other names: MDMB-4en-PINACA,
3 MDMB-PENINACA, and 5-CL-ADB-A.
- 4 [30] Methyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-
5 dimethyl-butanoate - Other names: 5F-MDMB-PICA and 5F-
6 MDMB-2201.
- 7 [31] 1-butyl-N-(1-carbamoyl-2,2-dimethyl-propyl)indazole-3-
8 carboxamide - Other names: ADB-BINACA and ADB-BUTINACA.
- 9 [32][30] 5-bromo-N-(1-carbamoyl-2,2-dimethyl-propyl)-1H-indazole-3-
10 carboxamide - Other names: ADB-5Br-INACA.
- 11 [33][31] Methyl 2-[(5-bromo-1H-indazole-3-carbonyl)amino]-3,3-dimethyl-
12 butanoate - Other names: MDMB-5Br-INACA.
- 13 [34][32] 5-bromo-1-butyl-N-(1-carbamoyl-2,2-dimethyl-propyl)indazole-3-
14 carboxamide - Other names: ADB-5'Br-BINACA and ADB-5'Br-
15 BUTINACA.
- 16 [33] Methyl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-
17 carboxamido)butanoate - Other name: MDMB-4en-PINACA.
- 18 [34] Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-
19 butanoate - Other names: 4F-MDMB-BUTICA; 4F-MDMB-BICA.
- 20 [35] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-en-1-yl)-1H-
21 indazole-3-carboxamide - Other name: ADB-4en-PINACA.
- 22 [36] Ethyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-dimethyl-
23 butanoate - Other names: 5F-EDMB-PICA; 5F-EDMB-2201.
- 24 [37] Methyl 2-(1-(4-fluorobenzyl)-1H-indole-3-carboxamido)-3-methyl
25 butanoate - Other name: MMB-FUBICA.
- 26 [38] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1H-indazole-3-
27 carboxamide - Other name: ADB-BUTINACA.
- 28 (4) Indole carboxylic acids. Any compound structurally derived from 1H-indole-
29 3-carboxylic acid or 1H-2-carboxylic acid substituted in both of the following
30 ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
31 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-

1 piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,
2 1-(N-methyl-3-morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo
3 benzyl group; and, at the hydroxyl group of the carboxylic acid by a phenyl,
4 benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, or propionaldehyde group
5 whether or not the compound is further modified to any extent in the
6 following ways:

- 7 (a) Substitution to the indole ring to any extent; or
8 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
9 cyclopropyl, propionaldehyde group to any extent; or
10 (c) A nitrogen heterocyclic analog of the indole ring; or
11 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,
12 adamantyl, or cyclopropyl ring.
13 (e) Examples include:

14 [1] 1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid 8-quinolinyl
15 ester - Other names: BB-22 and QUCHIC.

16 [2] naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate -
17 Other names: FDU-PB-22.

18 [3] 1-pentyl-1H-indole-3-carboxylic acid 8-quinolinyl ester - Other
19 names: PB-22 and QUPIC.

20 [4] 1-(5-Fluoropentyl)-1H-indole-3-carboxylic acid 8-quinolinyl ester -
21 Other names: 5-Fluoro PB-22 and 5F-PB-22.

22 [5] quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate - Other
23 names: FUB-PB-22.

24 [6] naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate -
25 Other names: NM2201 and CBL2201.

- 26 (5) Naphthylmethylindoles. Any compound containing a 1H-indol-3-yl-(1-
27 naphthyl)methane structure with substitution at the nitrogen atom of the
28 indole ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
29 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-
30 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or
31 (tetrahydropyran-4-yl)methyl group whether or not further substituted in the

- 1 indole ring to any extent and whether or not substituted in the naphthyl ring
2 to any extent. Examples include:
- 3 (a) 1-Pentyl-1H-indol-3-yl-(1-naphthyl)methane - Other names: JWH-175.
4 (b) 1-Pentyl-1H-indol-3-yl-(4-methyl-1-naphthyl)methane - Other names:
5 JWH-184.
- 6 (6) Naphthoylpyrroles. Any compound containing a 3-(1-naphthoyl)pyrrole
7 structure with substitution at the nitrogen atom of the pyrrole ring by an
8 alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
9 methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
10 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-
11 yl)methyl group whether or not further substituted in the pyrrole ring to any
12 extent, whether or not substituted in the naphthyl ring to any extent.
13 Examples include: (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-
14 ylmethanone - Other names: JWH-307.
- 15 (7) Naphthylmethylindenes. Any compound containing a naphthylideneindene
16 structure with substitution at the 3-position of the indene ring by an alkyl,
17 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-
18 2-piperidinyl)methyl, 2 (4 morpholinyl)ethyl, 1-(N-methyl-2-
19 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-
20 yl)methyl group whether or not further substituted in the indene ring to any
21 extent, whether or not substituted in the naphthyl ring to any extent.
22 Examples include: E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane
23 - Other names: JWH-176.
- 24 (8) Cyclohexylphenols. Any compound containing a 2-(3-
25 hydroxycyclohexyl)phenol structure with substitution at the 5-position of the
26 phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
27 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-
28 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or
29 (tetrahydropyran-4-yl)methyl group whether or not substituted in the
30 cyclohexyl ring to any extent. Examples include:

- 1 (a) 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
2 names: CP 47,497.
- 3 (b) 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
4 names: Cannabicyclohexanol and CP 47,497 C8 homologue.
- 5 (c) 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-
6 hydroxypropyl)cyclohexyl]-phenol - Other names: CP 55,940.
- 7 (9) Others specifically named:
- 8 (a) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-
9 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names: HU-210.
- 10 (b) (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-
11 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names:
12 Dexanabinol and HU-211.
- 13 (c) 2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-
14 benzoxazin-6-yl]-1-naphthalenylmethanone - Other names:
15 WIN 55,212-2.
- 16 (d) Naphthalen-1-yl-(4-pentyloxynaphthalen-1-yl)methanone - Other
17 names: CB-13.
- 18 (e) N-[(Z)-(1-hexyl-2-oxo-indolin-3-ylidene)amino]benzamide - Other
19 names: BZO-HEXOXIZID and MDA-19.
- 20 (f) N-[(Z)-(2-oxo-1-pentyl-indolin-3-ylidene)amino]benzamide - Other
21 names: BZO-POXIZID, Pentyl MDA-19, and 5C-MDA-19.
- 22 (g) N-[(Z)-[1-(5-fluoropentyl)-2-oxo-indolin-3-ylidene]amino]benzamide -
23 Other names: 5F-BZO-POXIZID and 5F-MDA-19.
- 24 (h) N-[(Z)-(2-oxo-1-pent-4-enyl-indolin-3-ylidene)amino]benzamide -
25 Other names: BZO-4en-POXIZID and 4en-pentyl MDA-19.
- 26 (i) N-[(Z)-[1-(cyclohexylmethyl)-2-oxo-indolin-3-ylidene]amino]benzamide
27 - Other names: BZO-CHMOXIZID, Cyclohexylmethyl MDA-19 and
28 CHM-MDA-19.
- 29 (j) N-(1-carbamoyl-2-methyl-propyl)-2-(5-fluoropentyl)-5-(4-
30 fluorophenyl)pyrazole-3-carboxamide - Other Names: 5F-AB-
31 PFUPPYCA.

- 1 (k) 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido[4,3-b]indol-1-one - Other
2 names: CUMYL-PEGACLONE; SGT-151.
- 3 o. Substituted phenethylamines. This includes any compound, unless specifically
4 excepted, specifically named in this schedule, or listed under a different
5 schedule, structurally derived from phenylethan-2-amine by substitution on the
6 phenyl ring in any of the following ways, that is to say, by substitution with a fused
7 methylenedioxy ring, fused furan ring, or fused tetrahydrofuran ring; by
8 substitution with two alkoxy groups; by substitution with one alkoxy and either
9 one fused furan, tetrahydrofuran, or tetrahydropyran ring system; or by
10 substitution with two fused ring systems from any combination of the furan,
11 tetrahydrofuran, or tetrahydropyran ring systems.
- 12 (1) Whether or not the compound is further modified in any of the following
13 ways, that is to say:
- 14 (a) By substitution of phenyl ring by any halo, hydroxyl, alkyl,
15 trifluoromethyl, alkoxy, or alkylthio groups;
- 16 (b) By substitution at the 2-position by any alkyl groups; or
- 17 (c) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl,
18 hydroxybenzyl, methylenedioxybenzyl, or methoxybenzyl groups.
- 19 (2) Examples include:
- 20 (a) 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (also known as 2C-C or
21 2,5-Dimethoxy-4-chlorophenethylamine).
- 22 (b) 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (also known as 2C-D or
23 2,5-Dimethoxy-4-methylphenethylamine).
- 24 (c) 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (also known as 2C-E or
25 2,5-Dimethoxy-4-ethylphenethylamine).
- 26 (d) 2-(2,5-Dimethoxyphenyl)ethanamine (also known as 2C-H or 2,5-
27 Dimethoxyphenethylamine).
- 28 (e) 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (also known as 2C-I or
29 2,5-Dimethoxy-4-iodophenethylamine).
- 30 (f) 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (also known as 2C-N or
31 2,5-Dimethoxy-4-nitrophenethylamine).

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

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

- 1 p. Substituted tryptamines. This includes any compound, unless specifically
2 excepted, specifically named in this schedule, or listed under a different
3 schedule, structurally derived from 2-(1H-indol-3-yl)ethanamine (i.e., tryptamine)
4 by mono- or di-substitution of the amine nitrogen with alkyl or alkenyl groups or
5 by inclusion of the amino nitrogen atom in a cyclic structure whether or not the
6 compound is further substituted at the alpha-position with an alkyl group or
7 whether or not further substituted on the indole ring to any extent with any alkyl,
8 alkoxy, halo, hydroxyl, or acetoxy groups. Examples include:
- 9 (1) 5-methoxy-N,N-diallyltryptamine (also known as 5-MeO-DALT).
10 (2) 4-acetoxy-N,N-dimethyltryptamine (also known as 4-AcO-DMT or O-
11 Acetylpsilocin).
12 (3) 4-hydroxy-N-methyl-N-ethyltryptamine (also known as 4-HO-MET).
13 (4) 4-hydroxy-N,N-diisopropyltryptamine (also known as 4-HO-DIPT).
14 (5) 5-methoxy-N-methyl-N-isopropyltryptamine (also known as 5-MeO-MiPT).
15 (6) 5-methoxy-N,N-dimethyltryptamine (also known as 5-MeO-DMT).
16 (7) Bufotenine (also known as 3-(Beta-Dimethyl-aminoethyl)-5-hydroxyindole;
17 3-(2-dimethylaminoethyl)-5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-
18 dimethyltryptamine; mappine).
19 (8) 5-methoxy-N,N-diisopropyltryptamine (also known as 5-MeO-DiPT).
20 (9) Diethyltryptamine (also known as N,N-Diethyltryptamine; DET).
21 (10) Dimethyltryptamine (also known as DMT).
22 (11) Psilocyn.
- 23 q. 1-[3-(trifluoromethylphenyl)]piperazine (also known as TFMPP).
24 r. 1-[4-(trifluoromethylphenyl)]piperazine.
25 s. 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (also known as 5,6-
26 Methylenedioxy-2-aminoindane or MDAI).
27 t. 2-(Ethylamino)-2-(3-methoxyphenyl)cyclohexanone (also known as
28 Methoxetamine or MXE).
29 u. Ethylamine analog of phencyclidine (also known as N-ethyl-1-
30 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl)
31 ethylamine, cyclohexamine, PCE).

- 1 v. Pyrrolidine analog of phencyclidine (also known as 1-(1-phenylcyclohexyl)-
2 pyrrolidine, PCPy, PHP).
- 3 w. Thiophene analog of phencyclidine (also known as (1-[1-(2-thienyl) cyclohexyl]
4 piperidine; 2-Thienylanalog of phencyclidine; TPCP, TCP).
- 5 x. 1-[1-(2-thienyl)cyclohexyl]pyrrolidine (also known as TCPy).
- 6 y. Salvia divinorum, salvinorin A, or any of the active ingredients of salvia divinorum.
- 7 6. Depressants. Unless specifically excepted or unless listed in another schedule, any
8 material compound, mixture, or preparation which contains any quantity of the
9 following substances having a depressant effect on the central nervous system,
10 whenever the existence of such salts, isomers, and salts of isomers is possible within
11 the specific chemical designation:
- 12 a. Gamma-hydroxybutyric acid.
- 13 b. Mecloqualone.
- 14 c. Methaqualone.
- 15 d. Clonazolam (also known as Clonitrazolam).
- 16 e. Etizolam.
- 17 f. Flualprazolam.
- 18 g. Flubromazepam.
- 19 h. Flubromazolam.
- 20 i. Adinazolam.
- 21 j. Bromazolam.
- 22 k. Deschloroetizolam.
- 23 l. Diclazepam.
- 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 (aa) N-ethylhexedrone (also known as alpha - ethylaminohexanophenone
20 and 2-(ethylamino)-1-phenylhexan-1-one)).
- 21 (bb) Alpha-pyrrolidinohexanophenone (also known as alpha-PHP, alpha-
22 pyrrolidinohexiophenone, and 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-
23 one)).
- 24 (cc) 4-methyl-alpha-ethylaminopentiophenone (also known as 4-MEAP
25 and 2-(ethylamino)-1-(4-methylphenyl)pentan-1-one)).
- 26 (dd) 4'-methyl-alpha-pyrrolidinohexiophenone (also known as MPHP, 4'-
27 methyl-alpha-pyrrolidinohexanophenone and 1-(4-methylphenyl)-2-
28 (pyrrolidin-1-yl)hexan-1-one)).
- 29 (ee) Alpha-pyrrolidinoheptaphenone (also known as PV8 and 1-phenyl-2-
30 (pyrrolidin-1-yl)heptan-1-one)).

- 1 (ff) 4-chloro-alpha-pyrrolidinovalerophenone (also known as 4-chloro-
2 alpha-PVP, 4'-chloro-alpha-pyrrolidinopentiophenone, and 1-(4-
3 chlorophenyl)-2-(pyrrolidin-1-yl)pentan-1-one)).
- 4 (gg) 4-methyl-1-phenyl-2-(pyrrolidin-1-yl)pentan-1-one (other name: alpha-
5 PiHP).
- 6 (hh) 2-(methylamino)-1-(3-methylphenyl)propan-1-one (other names: 3-
7 MMC; 3-methylmethcathinone).
- 8 (ii) Eutylone (also known as 1-(1,3-benzodioxol-5-yl)-2-
9 (ethylamino)butan-1-one).
- 10 d. Fenethylline.
- 11 e. Fluoroamphetamine.
- 12 f. Fluoromethamphetamine.
- 13 g. (±)cis-4-methylaminorex (also known as (±)cis-4,5-dihydro-4-methyl-5-phenyl-2-
14 oxazolamine).
- 15 h. N-Benzylpiperazine (also known as BZP, 1-benzylpiperazine).
- 16 i. N-ethylamphetamine.
- 17 j. N, N-dimethylamphetamine (also known as N,N-alpha-trimethyl-
18 benzeneethanamine; N,N-alpha-trimethylphenethylamine).
- 19 k. 1-(4-methoxyphenyl)-N-methylpropan-2-amine (also known as
20 paramethoxymethamphetamine and PMMA).
- 21 l. 4,4'-Dimethylaminorex (4,4'-DMAR; 4,5-dihydro-4-methyl-5-(4-methylphenyl)-2-
22 oxazolamine; 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine).
- 23 m. Amineptine (Also known as 7- [(10,11-dihydro-5Hdibenzo[a,d]cyclohepten-5-
24 yl)amino]heptanoic acid).
- 25 n. Mesocarb (Also known as N-phenyl-N'-(3-(1-phenylpropan-2-yl)-1,2,3-
26 oxadiazol-3-ium-5-yl)carbamimidate).
- 27 o. Methiopropamine (Also known as N-methyl-1-(thiophen-2-yl)propan-2-amine).
- 28 p. Ethylphenidate (ethyl 2-phenyl-2-(piperidin-2-yl)acetate).

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

1 **19-03.1-09. Schedule III.**

2 1. The controlled substances listed in this section are included in schedule III.

3 2. Schedule III consists of the drugs and other substances, by whatever official name,
4 common or usual name, chemical name, or brand name designated, listed in this
5 section.

6 3. Stimulants. Unless specifically excepted or unless listed in another schedule, any
7 material, compound, mixture, or preparation which contains any quantity of the
8 following substances having a stimulant effect on the central nervous system,
9 including its salts, isomers (whether optical, position, or geometric), and salts of such
10 isomers whenever the existence of such salts, isomers, and salts of isomers is
11 possible within the specific chemical designation:

12 a. Those compounds, mixtures, or preparations in dosage unit form containing any
13 stimulant substances listed in schedule II and any other drug of the quantitative
14 composition shown in that schedule for those drugs or which is the same except
15 that it contains a lesser quantity of controlled substances.

16 b. Benzphetamine.

17 c. Chlorphentermine.

18 d. Clortermine.

19 e. Phendimetrazine.

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

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

24 (1) Amobarbital;

25 (2) Secobarbital;

26 (3) Pentobarbital;

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

29 b. Any suppository dosage form containing:

30 (1) Amobarbital;

31 (2) Secobarbital;

- 1 (3) Pentobarbital;
2 or any salt of any of these drugs and approved by the food and drug
3 administration for marketing only as a suppository.
- 4 c. Any substance that contains any quantity of a derivative of barbituric acid, or any
5 salt of a derivative of barbituric acid, except those substances which are
6 specifically listed in other schedules thereof.
- 7 d. Chlorhexadol.
- 8 e. Embutramide.
- 9 f. Gamma-hydroxybutyric acid in a United States food and drug administration-
10 approved drug product.
- 11 g. Ketamine.
- 12 h. Lysergic acid.
- 13 i. Lysergic acid amide.
- 14 j. Methypylon.
- 15 k. Perampanel.
- 16 l. Sativex or its successor name as determined by the federal food and drug
17 administration.
- 18 m. Sulfondiethylmethane.
- 19 n. Sulfonethylmethane.
- 20 o. Sulfonmethane.
- 21 p. Tiletamine and zolazepam or any salt thereof. Some trade or other names for a
22 tiletamine-zolazepam combination product: Telazol. Some trade or other names
23 for tiletamine: 2-(ethylamino)-2-(2-thienyl)-cyclohexanone. Some trade or other
24 names for zolazepam: 4-2(2-fluorophenyl)-6, 8-dihydro-1,3,8-trimethylpyrazolo-
25 [3,4-e][1,4]-diazepin-7(1H)-one, flupyrazapon.
- 26 5. Nalorphine.
- 27 6. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any
28 material, compound, mixture, or preparation that contains any of the following narcotic
29 drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited
30 quantities as set forth below:

- 1 a. (1) Not more than 1.80 grams of codeine per 100 milliliters or not more than
2 90 milligrams per dosage unit, with an equal or greater quantity of an
3 isoquinoline alkaloid of opium.
- 4 (2) Not more than 1.80 grams of codeine per 100 milliliters or not more than
5 90 milligrams per dosage unit, with one or more active, non-narcotic
6 ingredients in recognized therapeutic amounts.
- 7 (3) Not more than 1.80 grams of dihydrocodeine per 100 milliliters or not more
8 than 90 milligrams per dosage unit, with one or more active, non-narcotic
9 ingredients in recognized therapeutic amounts.
- 10 (4) Not more than 300 milligrams of ethylmorphine per 100 milliliters or not
11 more than 15 milligrams per dosage unit, with one or more active,
12 non-narcotic ingredients in recognized therapeutic amounts.
- 13 (5) Not more than 500 milligrams of opium per 100 milliliters or per 100 grams,
14 or not more than 25 milligrams per dosage unit, with one or more active,
15 non-narcotic ingredients in recognized therapeutic amounts.
- 16 (6) Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams
17 with one or more active, non-narcotic ingredients in recognized therapeutic
18 amounts.
- 19 b. Buprenorphine.
- 20 7. Anabolic steroids. Unless specifically excepted or unless listed in another schedule,
21 any material, compound, mixture, or preparation that contains any quantity of the
22 following ~~anabolic steroid~~substances, including its salts, esters, and ethers:
- 23 a. 3beta,17-dihydroxy-5a-androstane;
- 24 b. 3alpha,17beta-dihydroxy-5a-androstane;
- 25 c. 5alpha-androstan-3,17-dione;
- 26 d. 5alpha-androstan-3,6,17-trione;
- 27 e. 1-androstenediol (3beta,17beta-dihydroxy-5alpha-androst-1-ene);
- 28 e.f. 1-androstenediol (3alpha,17beta-dihydroxy-5alpha-androst-1-ene);
- 29 f.g. 4-androstenediol (3beta,17beta-dihydroxyandrost-4-ene);
- 30 g.h. 5-androstenediol (3beta,17beta-dihydroxy-androst-5-ene);
- 31 h.i. 1-androstenedione ([5alpha]-androst-1-en-3,17-dione);

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- 1 i.j. 4-androstenedione (androst-4-en-3,17-dione);
- 2 j.k. 5-androstenedione (androst-5-en-3,17-dione);
- 3 k.l. Bolasterone (7alpha,17alpha-dimethyl-17beta-hydroxyandrost-4-en-3-one);
- 4 l.m. Boldenone (17beta-hydroxyandrost-1,4,-diene-3-one);
- 5 m.n. Boldione (androsta-1,4-diene-3,17-dione);
- 6 o. 6-bromo-androsta-1,4-diene-3,17-dione;
- 7 p. 6-bromo-androstan-3,17-dione;
- 8 n.q. Calusterone (7beta,17alpha-dimethyl-17beta-hydroxyandrost-4-en-3-one);
- 9 r. 4-chloro-17alpha-methyl-androsta-1,4-diene-3,17beta-diol;
- 10 s. 4-chloro-17alpha-methyl-androst-4-ene-3beta,17beta-diol;
- 11 t. 4-chloro-17alpha-methyl-17beta-hydroxy-androst-4-en-3-one;
- 12 u. 4-chloro-17alpha-methyl-17beta-hydroxy-androst-4-ene-3,11-dione;
- 13 e.v. Clostebol (4-chloro-17beta-hydroxyandrost-4-en-3-one);
- 14 p.w. Dehydrochloromethyltestosterone (4-chloro-17beta-hydroxy-17alpha-methyl-
15 androst-1,4-dien-3-one);
- 16 q.x. Delta-1-dihydrotestosterone (also known as '1-testosterone') (17beta-hydroxy-
17 5alpha-androst-1-en-3-one);
- 18 r.y. Desoxymethyltestosterone (17a-methyl-5a-androst-2-en-17beta-ol) (also known
19 as madol);
- 20 s.z. 4-dihydrotestosterone (17beta-hydroxy-androstan-3-one);
- 21 aa. 3beta,17beta-dihydroxy-5alpha-androstane;
- 22 bb. 3alpha,17beta-dihydroxy-5alpha-androstane;
- 23 cc. 2alpha,17alpha-dimethyl-17beta-hydroxy-5beta-androstan-3-one;
- 24 t.dd. Drostanolone (17beta-hydroxy-2alpha-methyl-5alpha-androstan-3-one);
- 25 ee. 2alpha,3alpha-epithio-17alpha-methyl-5alpha-androstan-17beta-ol;
- 26 ff. estra-4,9,11-triene-3,17-dione;
- 27 gg. 13beta-ethyl-17beta-hydroxygon-4-en-3-one;
- 28 u.hh. Ethylestrenol (17alpha-ethyl-17beta-hydroxyestr-4-ene);
- 29 v.ii. Fluoxymesterone (9-fluoro-17alpha-methyl-11beta, 17beta-dihydroxyandrost-4-
30 en-3-one);

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- 1 w.jj. Formebolone (2-formyl-17alpha-methyl-11alpha, 17beta-dihydroxyandrost-1,4-
2 dien-3-one);
- 3 x.kk. Furazabol (17alpha-methyl-17beta-hydroxyandrostando[2,3-c]-furazan);
- 4 ll. [3,2-c]furazan-5alpha-androstan-17beta-ol;
- 5 mm. 18a-homo-3-hydroxy-estra-2,5(10)-dien-17-one;
- 6 nn. 4-hydroxy-androst-4-ene-3,17-dione;
- 7 oo. 17beta-hydroxy-androstando[2,3-d]isoxazole;
- 8 pp. 17beta-hydroxy-androstando[3,2-c]isoxazole;
- 9 qq. 3beta-hydroxy-estra-4,9,11-trien-17-one;
- 10 y.rr. 13beta-ethyl-17alpha-hydroxygon-4-en-3-one;
- 11 z.ss. 4-hydroxytestosterone (4,17beta-dihydroxy-androst-4-en-3-one);
- 12 aa.tt. 4-hydroxy-19-nortestosterone (4,17beta-dihydroxy-estr-4-en-3-one);
- 13 bb.uu. Mestanolone (17alpha-methyl-17beta-hydroxy-5alpha-androstan-3-one);
- 14 ee.vv. Mesterolone (1alpha-methyl-17beta-hydroxy-[5alpha]-androstan-3-one);
- 15 dd.ww. Methandienone (17alpha-methyl-17beta-dihydroxyandrost-1,4-dien-3-one);
- 16 ee.xx. Methandriol (17alpha-methyl-3beta,17beta-dihydroxyandrost-5-ene);
- 17 ff.yy. Methasterone (2[alpha],17[alpha]-dimethyl-5[alpha]-androstan-17[beta]-ol-3-one);
- 18 gg.zz. Methenolone (1-methyl-17beta-hydroxy-5alpha-androst-1-en-3-one);
- 19 aaa. 17alpha-methyl-androsta-1,4-diene-3,17beta-diol;
- 20 bbb. 17alpha-methyl-5alpha-androstan-17beta-ol;
- 21 ccc. 17alpha-methyl-androstan-3-hydroxyimine-17beta-ol;
- 22 ddd. 6alpha-methyl-androst-4-ene-3,17-dione;
- 23 eee. 17alpha-methyl-androst-2-ene-3,17beta-diol;
- 24 hh.fff. 17alpha-methyl-3beta,17beta-dihydroxy-5a-androstane;
- 25 ii.ggg. 17alpha-methyl-3alpha,17beta-dihydroxy-5a-androstane;
- 26 jj.hhh. 17alpha-methyl-3beta,17beta-dihydroxyandrost-4-ene;
- 27 kk.iii. 17alpha-methyl-4-hydroxynandrolone (17alpha-methyl-4-hydroxy-17beta-
28 hydroxyestr-4-en-3-one);
- 29 H.jjj. Methyldienolone (17alpha-methyl-17beta-hydroxyestra-4,9(10)-dien-3-one);
- 30 mm.kkk. Methyltrienolone (17alpha-methyl-17beta-hydroxyestra-4,9(11)-trien-3-
31 one);

1 ~~###kkkk.~~ Trenbolone (17beta-hydroxyestr-4,9,11-trien-3-one);
2 ~~or any salt, ester, or isomer of a drug or substance described or listed in this~~
3 ~~subsection, if that salt, ester, or isomer promotes muscle growth.~~

4 The term does not include an anabolic steroid that is expressly intended for
5 administration through implants to cattle or other nonhuman species and which has
6 been approved by the secretary of health and human services for administration
7 unless any person prescribes, dispenses, possesses, delivers, or distributes for
8 human use.

9 8. Hallucinogenic substances.

10 a. Dronabinol (synthetic) [(-)-delta-9-(trans)-tetrahydrocannabinol] in sesame oil and
11 encapsulated in a soft gelatin capsule in a United States food and drug
12 administration-approved drug product.

13 b. Any product in hard or soft gelatin capsule form containing natural dronabinol
14 (derived from the cannabis plant) or synthetic dronabinol (produced from
15 synthetic materials) in sesame oil, for which an abbreviated new drug application
16 has been approved by the food and drug administration under section 505(j) of
17 the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355(j)] which references as
18 its listed drug the drug product referred to in subdivision a.

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

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

29 **19-03.1-11. Schedule IV.**

30 1. The controlled substances listed in this section are included in schedule IV.

- 1 2. Schedule IV consists of the drugs and other substances, by whatever official name,
2 common or usual name, chemical name, or brand name designated, listed in this
3 section.
- 4 3. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any
5 material, compound, mixture, or preparation containing any of the following narcotic
6 drugs or their salts calculated as the free anhydrous base or alkaloid, in limited
7 quantities as set forth below:
 - 8 a. Not more than 1 milligram of difenoxin and not less than 25 micrograms of
9 atropine sulfate per dosage unit.
 - 10 b. Dextropropoxyphene (also known as alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-
11 methyl-2-propionoxybutane).
 - 12 c. 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its salts, optical
13 and geometric isomers and salts of these isomers including Tramadol.
- 14 4. Depressants. Unless specifically excepted or unless listed in another schedule, any
15 material, compound, mixture, or preparation containing any quantity of the following
16 substances, including their salts, isomers, and salts of isomers whenever the
17 existence of those salts, isomers, and salts of isomers is possible within the specific
18 chemical designation:
 - 19 a. Alprazolam.
 - 20 b. Alfaxalone.
 - 21 c. Barbital.
 - 22 d. Brexanolone.
 - 23 e. Bromazepam.
 - 24 f. Camazepam.
 - 25 g. Carisoprodol.
 - 26 h. Chloral betaine.
 - 27 i. Chloral hydrate.
 - 28 j. Chlordiazepoxide.
 - 29 k. Clobazam.
 - 30 l. Clonazepam.
 - 31 m. Clorazepate.

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- 1 n. Clotiazepam.
- 2 o. Cloxazolam.
- 3 p. Daridorexant.
- 4 q. Delorazepam.
- 5 r. Diazepam.
- 6 s. Dichloralphenazone.
- 7 t. Estazolam.
- 8 u. Ethchlorvynol.
- 9 v. Ethinamate.
- 10 w. Ethyl loflazepate.
- 11 x. Fludiazepam.
- 12 y. Flunitrazepam.
- 13 z. Flurazepam.
- 14 aa. Fospropofol.
- 15 bb. Halazepam.
- 16 cc. Haloxazolam.
- 17 dd. Indiplon.
- 18 ee. Ketazolam.
- 19 ff. Lemborexant.
- 20 gg. Loprazolam.
- 21 hh. Lorazepam.
- 22 ii. Lorcaserin.
- 23 jj. Lormetazepam.
- 24 kk. Mebutamate.
- 25 ll. Medazepam.
- 26 mm. Meprobamate.
- 27 nn. Methohexital.
- 28 oo. Methylphenobarbital (also known as mephobarbital).
- 29 pp. Midazolam.
- 30 qq. Nimetazepam.
- 31 rr. Nitrazepam.

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- 1 ss. Nordiazepam.
- 2 tt. Oxazepam.
- 3 uu. Oxazolam.
- 4 vv. Paraldehyde.
- 5 ww. Petrichloral.
- 6 xx. Phenobarbital.
- 7 yy. Pinazepam.
- 8 zz. Propofol.
- 9 aaa. Prazepam.
- 10 bbb. Quazepam.
- 11 ccc. Remimazolam.
- 12 ddd. Suvorexant.
- 13 eee. Temazepam.
- 14 fff. Tetrazepam.
- 15 ggg. Triazolam.
- 16 hhh. Zaleplon.
- 17 iii. Zolpidem.
- 18 jjj. Zopiclone.
- 19 kkk. Zuranolone.
- 20 5. 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, and salts of isomers:
- 24 a. Cathine.
- 25 b. Diethylpropion.
- 26 c. Fencamfamin.
- 27 d. Fenproporex.
- 28 e. Mazindol.
- 29 f. Mefenorex.
- 30 g. Modafinil.
- 31 h. Pemoline (including organometallic complexes and chelates thereof).

- 1 i. Phentermine.
- 2 j. Pipradrol.
- 3 k. Serdexmethylphenidate.
- 4 l. Sibutramine.
- 5 m. Solriamfetol.
- 6 n. SPA ((-)-1-dimethylamino-1, 2-diphenylethane).
- 7 6. Other substances. Unless specifically excepted or unless listed in another schedule,
- 8 any material, compound, mixture, or preparation which contains any quantity of:
- 9 a. Pentazocine, including its salts.
- 10 b. Butorphanol, including its optical isomers.
- 11 c. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-
- 12 oxopropyl]][(1S)-1-(4-phenyl-1*H*-imidazol-2-yl)ethyl]amino]methyl]-2-
- 13 methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and
- 14 salts of isomers.
- 15 7. Hallucinogenic substances. Pharmaceutical composition of crystalline polymorph
- 16 psilocybin, known as COMP360 or any such trade name approved for COMP360 by
- 17 the United States food and drug administration.
- 18 8. The board may except by rule any compound, mixture, or preparation containing any
- 19 depressant substance listed in subsection 2 from the application of all or any part of
- 20 this chapter if the compound, mixture, or preparation contains one or more active
- 21 medicinal ingredients not having a depressant effect on the central nervous system,
- 22 and if the admixtures are included therein in combinations, quantity, proportion, or
- 23 concentration that vitiate the potential for abuse of the substances which have a
- 24 depressant effect on the central nervous system.

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