

Introduced by

Judiciary Committee

(At the request of the State Board of Pharmacy)

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

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

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

8 18. "Marijuana" means all parts of the plant *cannabis sativa* L., whether growing or not;
9 the seeds thereof; the resin extracted from any part of the plant; and every compound,
10 manufacture, salt, derivative, mixture, or preparation of the plant, its seeds, or resin.
11 The term does not include the mature stalks of the plant, fiber produced from the
12 stalks, oil or cake made from the seeds of the plant, any other compound,
13 manufacture, salt, derivative, mixture, or preparation of mature stalks, except the resin
14 extracted therefrom, fiber, oil, or cake, or the sterilized seed of the plant which is
15 incapable of germination. The term marijuana does not include hemp as defined in
16 title 4.1 chapter 4.1-18.1 or a prescription drug approved by the United States food and
17 drug administration under section 505 of the Federal Food, Drug, and Cosmetic Act
18 [21 U.S.C. 355].

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

21 **19-03.1-05. Schedule I.**

22 1. The controlled substances listed in this section are included in schedule I.

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

Sixty-seventh
Legislative Assembly

- 1 y. Hydroxypethidine.
- 2 z. Isotonitazene.
- 3 aa. Ketobemidone.
- 4 aa.bb. Levomoramide.
- 5 bb.cc. Levophenacymorphan.
- 6 ee.dd. Morpheridine.
- 7 dd.ee. MPPP (also known as 1-methyl-4-phenyl-4-propionoxypiperidine).
- 8 ee.ff. Noracymethadol.
- 9 ff.gg. Norlevorphanol.
- 10 gg.hh. Normethadone.
- 11 hh.ii. Norpipanone.
- 12 ii.jj. PEPAP (1-(2-Phenylethyl)-4-Phenyl-4-acetoxypiperidine).
- 13 jj.kk. Phenadoxone.
- 14 kk.ll. Phenampromide.
- 15 ll.mm. Phenomorphan.
- 16 mm.nn. Phenoperidine.
- 17 nn.oo. Piritramide.
- 18 oo.pp. Proheptazine.
- 19 pp.qq. Properidine.
- 20 qq.rr. Propiram.
- 21 rr.ss. Racemoramide.
- 22 ss.tt. Tilidine.
- 23 tt.uu. Trimeperidine.
- 24 uu.vv. 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as
- 25 U-47700).
- 26 vv.ww. 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine (also know as MT-45).
- 27 ww.xx. 3,4-dichloro-N-[[1-(dimethylamino)cyclohexyl]methyl]benzamide (also known as
- 28 AH-7921).
- 29 yy. Zipeprol.
- 30 xx.zz. Fentanyl derivatives. Unless specifically excepted or unless listed in another
- 31 schedule or are not FDA approved drugs, and are derived from N-(1-(2-

- 1 Phenylethyl)-4-piperidiny]-N-phenylpropanamide (Fentanyl) by any substitution
2 on or replacement of the phenethyl group, any substitution on the piperidine ring,
3 any substitution on or replacement of the propanamide group, any substitution on
4 the anilido phenyl group, or any combination of the above. Examples include:
- 5 (1) N-[1-(1-methyl-2-phenethyl)-4-piperidiny]-N-phenylacetamide (also known
6 as Acetyl-alpha-methylfentanyl).
 - 7 (2) N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-
8 2-phenylethyl)-4-(N-propanilido)piperidine (also known as Alpha-
9 methylfentanyl).
 - 10 (3) N-[1-methyl-2-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide (also
11 known as Alpha-methylthiofentanyl).
 - 12 (4) N-[1-(2-hydroxy-2-phenethyl)-4-piperidiny]-N-phenylpropanamide (also
13 known as Beta-hydroxyfentanyl).
 - 14 (5) N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidiny]-N-phenylpropanamide
15 (also known as Beta-hydroxy-3-methylfentanyl).
 - 16 (6) N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (also
17 known as 3-Methylfentanyl).
 - 18 (7) N-[3-methyl-1-(2-thienyl)ethyl-4-piperidiny]-N-phenylpropanamide (also
19 known as 3-Methylthiofentanyl).
 - 20 (8) N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidiny]propanamide (also
21 known as Para-fluorofentanyl).
 - 22 (9) N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidiny]propanamide (also known as
23 Thiofentanyl).
 - 24 (10) N-(1-phenylethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide (also known
25 as Furanyl Fentanyl).
 - 26 (11) N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide; N-(1-phenethylpiperidin-
27 4-yl)-N-phenylbutanamide (also known as Butyryl Fentanyl).
 - 28 (12) N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide;
29 N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidiny]-N-phenylpropanamide (also
30 known as Beta-Hydroxythiofentanyl).

Sixty-seventh
Legislative Assembly

- 1 (13) N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as Acetyl
2 Fentanyl).
- 3 (14) N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (also known as Acryl
4 Fentanyl).
- 5 (15) N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide (also known as Valeryl
6 Fentanyl).
- 7 (16) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known
8 as 4-Fluoroisobutyryl Fentanyl).
- 9 (17) N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide (also known
10 as Ortho-fluorofentanyl, 2-Fluorofentanyl).
- 11 (18) N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (also
12 known as Tetrahydrofuranyl Fentanyl).
- 13 (19) 2-methoxy-N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (also known as
14 Methoxyacetyl Fentanyl).
- 15 (20) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (also
16 known as Cyclopropyl Fentanyl).
- 17 (21) N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetamide (also
18 known as Ocfentanil).
- 19 (22) N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (also
20 known as Cyclopentyl Fentanyl).
- 21 (23) N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (also known as
22 Isobutyryl Fentanyl).
- 23 (24) N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known
24 as Para-chloroisobutyryl Fentanyl).
- 25 (25) N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known
26 as Para-methoxybutyryl Fentanyl).
- 27 (26) N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known as
28 Para-fluorobutyryl Fentanyl).
- 29 (27) N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide (also
30 known as 2'-fluoro Ortho-fluorofentanyl).

1 (28) N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide (also known as
2 Ortho-methyl Acetylfentanyl).

3 (29) N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide (also known as
4 Beta'-phenyl Fentanyl and Hydrocinnamoyl Fentanyl).

5 (30) N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide (also
6 known as Thiofuranyl Fentanyl).

7 (31) (E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide (also known as
8 Crotonyl Fentanyl).

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

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

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

- 1 n. Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained in a
2 plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of
3 the substances contained in the cannabis plant, or in the resinous extractives of
4 such plant, including synthetic substances, derivatives, and their isomers with
5 similar chemical structure and pharmacological activity to those substances
6 contained in the plant; excluding tetrahydrocannabinols found in hemp as defined
7 in title 4.1; such as the following:
- 8 (1) Delta-1 cis or trans tetrahydrocannabinol, and their optical isomers. Other
9 names: Delta-9-tetrahydrocannabinol.
- 10 (2) Delta-6 cis or trans tetrahydrocannabinol, and their optical isomers.
- 11 (3) Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers.
12 (Since nomenclature of these substances is not internationally standardized,
13 compounds of these structures, regardless of numerical designation of atomic
14 positions covered.)
- 15 o. Cannabinoids, synthetic. It includes the chemicals and chemical groups listed
16 below, including their homologues, salts, isomers, and salts of isomers. The term
17 "isomer" includes the optical, position, and geometric isomers.
- 18 (1) Indole carboxaldehydes. Any compound structurally derived from 1H-indole-
19 3-carboxaldehyde or 1H-2-carboxaldehyde substituted in both of the
20 following ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
21 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
22 piperidiny)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidiny)methyl,
23 1-(N-methyl-3- morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo
24 benzyl group; and, at the hydrogen of the carboxaldehyde by a phenyl,
25 benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, pyrrolidiny, piperaziny, or
26 propionaldehyde group whether or not the compound is further modified to
27 any extent in the following ways:
- 28 (a) Substitution to the indole ring to any extent; or
29 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
30 cyclopropyl, pyrrolidiny, piperaziny, or propionaldehyde group to any
31 extent; or

- 1 (c) A nitrogen heterocyclic analog of the indole ring; or
- 2 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,
- 3 adamantyl, or cyclopropyl ring.
- 4 (e) Examples include:
- 5 [1] 1-Pentyl-3-(1-naphthoyl)indole - Other names: JWH-018 and
- 6 AM-678.
- 7 [2] 1-Butyl-3-(1-naphthoyl)indole - Other names: JWH-073.
- 8 [3] 1-Pentyl-3-(4-methoxy-1-naphthoyl)indole - Other names:
- 9 JWH-081.
- 10 [4] 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole - Other names:
- 11 JWH-200.
- 12 [5] 1-Propyl-2-methyl-3-(1-naphthoyl)indole - Other names:
- 13 JWH-015.
- 14 [6] 1-Hexyl-3-(1-naphthoyl)indole - Other names: JWH-019.
- 15 [7] 1-Pentyl-3-(4-methyl-1-naphthoyl)indole - Other names:
- 16 JWH-122.
- 17 [8] 1-Pentyl-3-(4-ethyl-1-naphthoyl)indole - Other names: JWH-210.
- 18 [9] 1-Pentyl-3-(4-chloro-1-naphthoyl)indole - Other names:
- 19 JWH-398.
- 20 [10] 1-(5-fluoropentyl)-3-(1-naphthoyl)indole - Other names:
- 21 AM-2201.
- 22 [11] 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole - Other
- 23 names: RCS-8.
- 24 [12] 1-Pentyl-3-(2-methoxyphenylacetyl)indole - Other names:
- 25 JWH-250.
- 26 [13] 1-Pentyl-3-(2-methylphenylacetyl)indole - Other names:
- 27 JWH-251.
- 28 [14] 1-Pentyl-3-(2-chlorophenylacetyl)indole - Other names: JWH-
- 29 203.
- 30 [15] 1-Pentyl-3-(4-methoxybenzoyl)indole - Other names: RCS-4.

- 1 [16] (1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole) - Other names:
2 AM-694.
- 3 [17] (4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-
4 yl]methanone - Other names: WIN 48,098 and Pravadoline.
- 5 [18] (1-Pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone --
6 Other names: UR-144.
- 7 [19] (1-(5-fluoropentyl)indol-3-yl)-(2,2,3,3-
8 tetramethylcyclopropyl)methanone - Other names: XLR-11.
- 9 [20] (1-(2-morpholin-4-ylethyl)-1H-indol-3-yl)-(2,2,3,3-
10 tetramethylcyclopropyl)methanone - Other names: A-796,260.
- 11 [21] (1-(5-fluoropentyl)-1H-indazol-3-yl)(naphthalen-1-yl)methanone --
12 Other names: THJ-2201.
- 13 [22] 1-naphthalenyl(1-pentyl-1H-indazol-3-yl)-methanone -- Other
14 names: THJ-018.
- 15 [23] (1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-yl)(naphthalen-1-
16 yl)methanone - Other names: FUBIMINA.
- 17 [24] 1-[(N-methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl) indole -
18 Other names: AM-1248.
- 19 [25] 1-Pentyl-3-(1-adamantoyl)indole - Other names: AB-001 and
20 JWH-018 adamantyl analog.
- 21 (2) Indole carboxamides. Any compound structurally derived from 1H-indole-3-
22 carboxamide or 1H-2-carboxamide substituted in both of the following ways:
23 at the nitrogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,
24 alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,
25 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-
26 morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo benzyl group;
27 and, at the nitrogen of the carboxamide by a phenyl, benzyl, cumyl,
28 naphthyl, adamantyl, cyclopropyl, or propionaldehyde group whether or not
29 the compound is further modified to any extent in the following ways:
30 (a) Substitution to the indole ring to any extent; or

- 1 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
2 cyclopropyl, or propionaldehyde group to any extent; or
3 (c) A nitrogen heterocyclic analog of the indole ring; or
4 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,
5 adamantyl, or cyclopropyl ring.
6 (e) Examples include:
7 [1] N-Adamantyl-1-pentyl-1H-indole-3-carboxamide - Other names:
8 JWH-018 adamantyl carboxamide, APICA, SDB-001, and 2NE1.
9 [2] N-Adamantyl-1-fluoropentylindole-3-carboxamide - Other names:
10 STS-135.
11 [3] N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide - Other
12 names: AKB 48 and APINACA.
13 [4] N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide - Other
14 names: NNEI and MN-24.
15 [5] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-
16 carboxamide - Other names: ADBICA.
17 [6] (S)-N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-
18 3-carboxamide - Other names: AB-PINACA.
19 [7] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-
20 fluorophenyl)methyl]-1H-indazole-3-carboxamide - Other names:
21 AB-FUBINACA.
22 [8] N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-
23 indazole-3-carboxamide - Other names: 5-Fluoro AB-PINACA
24 and 5F-AB-PINACA.
25 [9] N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-
26 3-carboxamide - Other names: ADB-PINACA.
27 [10] N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-(cyclohexylmethyl)-
28 1H-indazole-3-carboxamide - Other names: AB-CHMINACA.
29 [11] N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-
30 indazole-3-carboxamide - Other names: ADB-FUBINACA.

- 1 [12] ~~N-((3s,5s,7s)-adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-~~
2 ~~carboxamide - Other names: FUB-AKB48 and AKB48 N-(4-~~
3 ~~fluorobenzyl) analog~~N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H
4 ~~-indazole-3-carboxamide - Other names: FUB-AKB48, FUB-~~
5 ~~APINACA, and 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.
- 10 [26] 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-
11 carboxamide - Other names: 5F-CUMYL-PINACA, SGT-25).
- 12 [27] (1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)
13 methanone - Other names: FUB-144.
- 14 (3) Indole carboxylic acids. Any compound structurally derived from 1H-indole-
15 3-carboxylic acid or 1H-2-carboxylic acid substituted in both of the following
16 ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl,
17 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
18 piperidiny)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidiny)methyl,
19 1-(N-methyl-3-morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo
20 benzyl group; and, at the hydroxyl group of the carboxylic acid by a phenyl,
21 benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, or propionaldehyde group
22 whether or not the compound is further modified to any extent in the
23 following ways:
- 24 (a) Substitution to the indole ring to any extent; or
- 25 (b) Substitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
26 cyclopropyl, propionaldehyde group to any extent; or
- 27 (c) A nitrogen heterocyclic analog of the indole ring; or
- 28 (d) A nitrogen heterocyclic analog of the phenyl, benzyl, naphthyl,
29 adamantyl, or cyclopropyl ring.
- 30 (e) Examples include:

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

- 1 Examples include: (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-
2 ylmethanone - Other names: JWH-307.
- 3 (6) Naphthylmethylindenes. Any compound containing a naphthylideneindene
4 structure with substitution at the 3-position of the indene ring by an alkyl,
5 haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-
6 2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
7 pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-
8 yl)methyl group whether or not further substituted in the indene ring to any
9 extent, whether or not substituted in the naphthyl ring to any extent.
- 10 Examples include: E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane
11 - Other names: JWH-176.
- 12 (7) Cyclohexylphenols. Any compound containing a 2-(3-
13 hydroxycyclohexyl)phenol structure with substitution at the 5-position of the
14 phenolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
15 cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-
16 (N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or
17 (tetrahydropyran-4-yl)methyl group whether or not substituted in the
18 cyclohexyl ring to any extent. Examples include:
- 19 (a) 5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
20 names: CP 47,497.
- 21 (b) 5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
22 names: Cannabicyclohexanol and CP 47,497 C8 homologue.
- 23 (c) 5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-
24 hydroxypropyl)cyclohexyl]-phenol - Other names: CP 55,940.
- 25 (8) Others specifically named:
- 26 (a) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-
27 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names: HU-210.
- 28 (b) (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-
29 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names:
30 Dexanabinol and HU-211.

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

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

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

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

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

- 1 a. Aminorex (also known as 2-amino-5-phenyl-2-oxazoline, or 4,5-dihydro-5-phenyl-
2 2-oxazolamine).
- 3 b. Cathinone.
- 4 c. Substituted cathinones. Any compound, material, mixture, preparation, or other
5 product, unless listed in another schedule or an approved food and drug
6 administration drug (e.g., bupropion, pyrovalerone), structurally derived from 2-
7 aminopropan-1-one by substitution at the 1-position with either phenyl, naphthyl,
8 or thiophene ring systems, whether or not the compound is further modified in
9 any of the following ways:
- 10 (1) By substitution in the ring system to any extent with alkyl, alkylendioxy,
11 alkoxy, haloalkyl, hydroxyl, or halide substituents, whether or not further
12 substituted in the ring system by one or more other univalent substituents;
- 13 (2) By substitution at the 3-position with an acyclic alkyl substituent;
- 14 (3) By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or
15 methoxybenzyl groups; or
- 16 (4) By inclusion of the 2-amino nitrogen atom in a cyclic structure.
- 17 Some trade or other names:
- 18 (a) 3,4-Methylenedioxy-alpha-pyrrolidinopropiophenone (also known as
19 MDPPP).
- 20 (b) 3,4-Methylenedioxy-N-ethylcathinone (also known as Ethylone,
21 MDEC, or bk-MDEA).
- 22 (c) 3,4-Methylenedioxy-N-methylcathinone (also known as Methylone or
23 bk-MDMA).
- 24 (d) 3,4-Methylenedioxyprovalerone (also known as MDPV).
- 25 (e) 3,4-Dimethylmethcathinone (also known as 3,4-DMMC).
- 26 (f) 2-(methylamino)-1-phenylpentan-1-one (also known as Pentedrone).
- 27 (g) 2-Fluoromethcathinone (also known as 2-FMC).
- 28 (h) 3-Fluoromethcathinone (also known as 3-FMC).
- 29 (i) 4-Methylethcathinone (also known as 4-MEC and 4-methyl-N-
30 ethylcathinone).
- 31 (j) 4-Fluoromethcathinone (also known as Flephedrone and 4-FMC).

Sixty-seventh
Legislative Assembly

- 1 (k) 4-Methoxy-alpha-pyrrolidinopropiophenone (also known as MOPPP).
2 (l) 4-Methoxymethcathinone (also known as Methedrone; bk-PMMA).
3 (m) 4'-Methyl-alpha-pyrrolidinobutiophenone (also known as MPBP).
4 (n) Alpha-methylamino-butyrophenone (also known as Buphedrone or
5 MABP).
6 (o) Alpha-pyrrolidinobutiophenone (also known as alpha-PBP).
7 (p) Alpha-pyrrolidinopropiophenone (also known as alpha-PPP).
8 (q) Alpha-pyrrolidinopentiophenone (also known as Alpha-
9 pyrrolidinovalerophenone or alpha-PVP).
10 (r) Beta-keto-N-methylbenzodioxolylbutanamine (also known as Butylone
11 or bk-MBDB).
12 (s) Ethcathinone (also known as N-Ethylcathinone).
13 (t) 4-Methylmethcathinone (also known as Mephedrone or 4-MMC).
14 (u) Methcathinone.
15 (v) N,N-dimethylcathinone (also known as metamfepramone).
16 (w) Naphthylpyrovalerone (naphyrone).
17 (x) B-Keto-Methylbenzodioxolylpentanamine (also known as Pentylone).
18 (y) 4-Methyl-alpha-pyrrolidinopropiophenone (also known as 4-MePPP
19 and MPPP).
20 (z) 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one (also known as
21 Ephylone and N-Ethylpentylone).
22 (aa) N-ethylhexedrone.
23 (bb) alpha-pyrrolidinohexanophenone (also known as alpha-PHP).
24 (cc) 4-methyl-alpha-ethylaminopentiophenone (also known as 4-MEAP).
25 (dd) 4'-methyl-alpha-pyrrolidinohexiophenone (also known as MPHP).
26 (ee) alpha-pyrrolidinoheptaphenone (also known as PV8).
27 (ff) 4-chloro-alpha-pyrrolidinovalerophenone (also known 4-chloro-alpha-
28 PVP).
29 d. Fenethylline.
30 e. Fluoroamphetamine.
31 f. Fluoromethamphetamine.

- 1 g. (±)cis-4-methylaminorex (also known as (±)cis-4,5-dihydro-4-methyl-5-phenyl-2-
- 2 oxazolamine).
- 3 h. N-Benzylpiperazine (also known as BZP, 1-benzylpiperazine).
- 4 i. N-ethylamphetamine.
- 5 j. N, N-dimethylamphetamine (also known as N,N-alpha-trimethyl-
- 6 benzeneethanamine; N,N-alpha-trimethylphenethylamine).

7 **SECTION 3. AMENDMENT.** Section 19-03.1-07 of the North Dakota Century Code is
8 amended and reenacted as follows:

9 **19-03.1-07. Schedule II.**

- 10 1. The controlled substances listed in this section are included in schedule II.
- 11 2. Schedule II consists of the drugs and other substances, by whatever official name,
12 common or usual name, chemical name, or brand name designated, listed in this
13 section.
- 14 3. Substances, vegetable origin or chemical synthesis. Unless specifically excepted or
15 unless listed in another schedule, any of the following substances whether produced
16 directly or indirectly by extraction from substances of vegetable origin, or
17 independently by means of chemical synthesis, or by a combination of extraction and
18 chemical synthesis:
 - 19 a. Opium and opiate, and any salt, compound, derivative, or preparation of opium or
20 opiate, excluding apomorphine, thebaine-derived butorphanol, dextrophan,
21 nalbuphine, naldemedine, nalmefene, naloxegol, naloxone, 6 beta-naltrexol, and
22 naltrexone and their respective salts, but including the following:
 - 23 (1) Codeine.
 - 24 (2) Dihydroetorphine.
 - 25 (3) Ethylmorphine.
 - 26 (4) Etorphine hydrochloride.
 - 27 (5) Granulated opium.
 - 28 (6) Hydrocodone.
 - 29 (7) Hydromorphone.
 - 30 (8) Metopon.
 - 31 (9) Morphine.

- 1 (10) Noroxymorphone.
- 2 (11) Opium extracts.
- 3 ~~(11)~~~~(12)~~ Opium fluid.
- 4 ~~(12)~~~~(13)~~ Oripavine.
- 5 ~~(13)~~~~(14)~~ Oxycodone.
- 6 ~~(14)~~~~(15)~~ Oxymorphone.
- 7 ~~(15)~~~~(16)~~ Powder opium.
- 8 ~~(16)~~~~(17)~~ Raw opium.
- 9 ~~(17)~~~~(18)~~ Thebaine.
- 10 ~~(18)~~~~(19)~~ Tincture of opium.
- 11 b. Any salt, compound, derivative, or preparation thereof which is chemically
- 12 equivalent or identical with any of the substances referred to in subdivision a, but
- 13 not including the isoquinoline alkaloids of opium.
- 14 c. Opium poppy and poppy straw.
- 15 d. Coca leaves and any salt, compound, derivative, or preparation of coca leaves,
- 16 including cocaine and ecgonine and their salts, isomers, derivatives, and salts of
- 17 isomers and derivatives, and any salt, compound, derivative, or preparation
- 18 thereof that is chemically equivalent or identical with any of these substances,
- 19 except that the nondosage substances must include decocainized coca leaves or
- 20 extractions of coca leaves which do not contain cocaine or ecgonine.
- 21 e. Concentrate of poppy straw (the crude extract of poppy straw in either liquid,
- 22 solid, or powder form which contains the phenanthrine alkaloids of the opium
- 23 poppy).
- 24 4. Opiates. Unless specifically excepted or unless in another schedule, any of the
- 25 following opiates, including their isomers, esters, ethers, salts, and salts of isomers,
- 26 esters, and ethers whenever the existence of those isomers, esters, ethers, and salts
- 27 is possible within the specific chemical designation, dextrophan and
- 28 levopropoxyphene excepted:
- 29 a. Alfentanil.
- 30 b. Alphaprodine.
- 31 c. Anileridine.

Sixty-seventh
Legislative Assembly

- 1 d. Bezitramide.
- 2 e. Bulk dextropropoxyphene (nondosage forms).
- 3 f. Carfentanil.
- 4 g. Dihydrocodeine.
- 5 h. Diphenoxylate.
- 6 i. Fentanyl.
- 7 j. Isomethadone.
- 8 k. Levo-alphaacetylmethadol (LAAM).
- 9 l. Levomethorphan.
- 10 m. Levorphanol.
- 11 n. Metazocine.
- 12 o. Methadone.
- 13 p. Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane.
- 14 q. Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic
15 acid.
- 16 r. Oliceridine (N-[(3-methoxythiophen-2-yl)methyl] (2-[(9R)-9-(pyridin-2-yl)-6-
17 oxaspiro [4.5]decan-9-yl]ethyl))amine fumarate).
- 18 s. Pethidine (also known as meperidine).
- 19 s.t. Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine.
- 20 t.u. Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate.
- 21 u.v. Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid.
- 22 v.w. Phenazocine.
- 23 w.x. Priminodine.
- 24 x.y. Racemethorphan.
- 25 y.z. Racemorphan.
- 26 z.aa. Remifentanil.
- 27 aa.bb. Sufentanil.
- 28 bb.cc. Tapentadol.
- 29 ee.dd. Thiafentanil.

- 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 a. Amphetamine, its salts, optical isomers, and salts of its optical isomers.
5 b. Lisdexamfetamine, its salts, isomers, and salts of isomers.
6 c. Methamphetamine, its salts, isomers, and salts of isomers.
7 d. Phenmetrazine and its salts.
8 e. Methylphenidate.
- 9 6. Depressants. 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 depressant effect on the central nervous system,
12 including its salts, isomers, and salts of isomers whenever the existence of such salts,
13 isomers, and salts of isomers is possible within the specific chemical designation:
14 a. Amobarbital.
15 b. Glutethimide.
16 c. Pentobarbital.
17 d. Phencyclidine.
18 e. Secobarbital.
- 19 7. Hallucinogenic substances.
20 a. Nabilone [another name for nabilone (\pm)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8,
21 10, 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9Hdibenzo [b, d] pyran-9-one].
22 b. Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] in an oral solution in a drug
23 product approved for marketing by the federal food and drug administration.
- 24 8. Immediate precursors. Unless specifically excepted or unless listed in another
25 schedule, any material, compound, mixture, or preparation that contains any quantity
26 of the following substances:
27 a. Immediate precursor to amphetamine and methamphetamine: Phenylacetone.
28 Some trade or other names: phenyl-2-propanone; P2P, benzyl methyl ketone;
29 methyl benzyl ketone.
30 b. Immediate precursors to phencyclidine (PCP):
31 (1) 1-phenylcyclohexylamine.

1 (2) 1-piperidinocyclohexanecarbonitrile (PCC).

2 c. Immediate precursors to fentanyl: 4-anilino-N-phenethyl-4-piperidine (ANPP).

3 **SECTION 4. AMENDMENT.** Section 19-03.1-11 of the North Dakota Century Code is
4 amended and reenacted as follows:

5 **19-03.1-11. Schedule IV.**

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

7 2. Schedule IV consists of the drugs and other substances, by whatever official name,
8 common or usual name, chemical name, or brand name designated, listed in this
9 section.

10 3. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any
11 material, compound, mixture, or preparation containing 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. Not more than 1 milligram of difenoxin and not less than 25 micrograms of
15 atropine sulfate per dosage unit.

16 b. Dextropropoxyphene (also known as alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-
17 methyl-2-propionoxybutane).

18 c. 2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its salts, optical
19 and geometric isomers and salts of these isomers including Tramadol.

20 4. Depressants. Unless specifically excepted or unless listed in another schedule, any
21 material, compound, mixture, or preparation containing any quantity of the following
22 substances, including their salts, isomers, and salts of isomers whenever the
23 existence of those salts, isomers, and salts of isomers is possible within the specific
24 chemical designation:

25 a. Alprazolam.

26 b. Alfaxalone.

27 c. Barbital.

28 d. Brexanolone.

29 e. Bromazepam.

30 e-f. Camazepam.

31 f-g. Carisoprodol.

Sixty-seventh
Legislative Assembly

1	<u>g-h.</u>	Chloral betaine.
2	<u>h-i.</u>	Chloral hydrate.
3	<u>i-j.</u>	Chlordiazepoxide.
4	<u>j-k.</u>	Clobazam.
5	<u>k-l.</u>	Clonazepam.
6	<u>l-m.</u>	Clorazepate.
7	<u>m-n.</u>	Clotiazepam.
8	<u>n-o.</u>	Cloxazolam.
9	<u>o-p.</u>	Delorazepam.
10	<u>p-q.</u>	Diazepam.
11	<u>q-r.</u>	Dichloralphenazone.
12	<u>r-s.</u>	Estazolam.
13	<u>s-t.</u>	Ethchlorvynol.
14	<u>t-u.</u>	Ethinamate.
15	<u>u-v.</u>	Ethyl loflazepate.
16	<u>v-w.</u>	Fludiazepam.
17	<u>w-x.</u>	Flunitrazepam.
18	<u>x-y.</u>	Flurazepam.
19	<u>y-z.</u>	Fospropofol.
20	<u>z-aa.</u>	Halazepam.
21	<u>aa-bb.</u>	Haloxazolam.
22	<u>bb-cc.</u>	Indiplon.
23	<u>cc-dd.</u>	Ketazolam.
24	<u>dd-ee.</u>	<u>Lemborexant.</u>
25	<u>ff.</u>	Loprazolam.
26	<u>ee-gg.</u>	Lorazepam.
27	<u>ff-hh.</u>	Lorcaserin.
28	<u>gg-ii.</u>	Lormetazepam.
29	<u>hh-jj.</u>	Mebutamate.
30	<u>ii-kk.</u>	Medazepam.
31	<u>jj-ll.</u>	Meprobamate.

Sixty-seventh
Legislative Assembly

- 1 kk.mm. Methohexital.
- 2 ll.nn. Methylphenobarbital (also known as mephobarbital).
- 3 mm.oo. Midazolam.
- 4 nn.pp. Nimetazepam.
- 5 oo.qq. Nitrazepam.
- 6 pp.rr. Nordiazepam.
- 7 qq.ss. Oxazepam.
- 8 rr.tt. Oxazolam.
- 9 ss.uu. Paraldehyde.
- 10 tt.vv. Petrichloral.
- 11 uu.wv. Phenobarbital.
- 12 vv.xx. Pinazepam.
- 13 ww.yy. Propofol.
- 14 xx.zz. Prazepam.
- 15 yy.aaa. Quazepam.
- 16 zz.bbb. Remimazolam.
- 17 ccc. Suvorexant.
- 18 aaa.ddd. Temazepam.
- 19 bbb.eee. Tetrazepam.
- 20 eee.fff. Triazolam.
- 21 ddd.ggg. Zaleplon.
- 22 eee.hhh. Zolpidem.
- 23 fff.iii. Zopiclone.
- 24 5. Fenfluramine. Any material, compound, mixture, or preparation which contains any
- 25 quantity of the following substances, including its salts, isomers (whether optical,
- 26 position, or geometric), and salts of such isomers, whenever the existence of such
- 27 salts, isomers, and salts of isomers is possible: Fenfluramine.
- 28 6. Stimulants. Unless specifically excepted or unless listed in another schedule, any
- 29 material, compound, mixture, or preparation which contains any quantity of the
- 30 following substances having a stimulant effect on the central nervous system,
- 31 including its salts, isomers, and salts of isomers:

Sixty-seventh
Legislative Assembly

- 1 a. Cathine.
- 2 b. Diethylpropion.
- 3 c. Fencamfamin.
- 4 d. Fenproporex.
- 5 e. Mazindol.
- 6 f. Mefenorex.
- 7 g. Modafinil.
- 8 h. Pemoline (including organometallic complexes and chelates thereof).
- 9 i. Phentermine.
- 10 j. Pipradrol.
- 11 k. Sibutramine.
- 12 l. Solriamfetol.
- 13 m. SPA ((-)-1-dimethylamino-1, 2-diphenylethane).
- 14 7. Other substances. Unless specifically excepted or unless listed in another schedule,
15 any material, compound, mixture, or preparation which contains any quantity of:
 - 16 a. Pentazocine, including its salts.
 - 17 b. Butorphanol, including its optical isomers.
 - 18 c. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl]-2,6-dimethylphenyl]-1-
19 oxopropyl]][(1S)-1-(4-phenyl-1*H*-imidazol-2-yl)ethyl]amino]methyl]-2-
20 methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and
21 salts of isomers.
- 22 8. The board may except by rule any compound, mixture, or preparation containing any
23 depressant substance listed in subsection 2 from the application of all or any part of
24 this chapter if the compound, mixture, or preparation contains one or more active
25 medicinal ingredients not having a depressant effect on the central nervous system,
26 and if the admixtures are included therein in combinations, quantity, proportion, or
27 concentration that vitiate the potential for abuse of the substances which have a
28 depressant effect on the central nervous system.

29 **SECTION 5. AMENDMENT.** Section 19-03.1-13 of the North Dakota Century Code is
30 amended and reenacted as follows:

1 **19-03.1-13. Schedule V.**

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

3 2. Schedule V 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. Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any
7 material, compound, mixture, or preparation containing any of the following narcotic
8 drugs and their salts.

9 4. Narcotic drugs containing non-narcotic active medicinal ingredients. Any compound,
10 mixture, or preparation containing any of the following narcotic drugs, or their salts
11 calculated as the free anhydrous base or alkaloid, in limited quantities as set forth
12 below, which includes one or more non-narcotic active medicinal ingredients in
13 sufficient proportion to confer upon the compound, mixture, or preparation valuable
14 medicinal qualities other than those possessed by narcotic drugs alone.

15 a. Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams.

16 b. Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per
17 100 grams.

18 c. Not more than 100 milligrams of ethylmorphine per 100 milliliters or per
19 100 grams.

20 d. Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms
21 of atropine sulfate per dosage unit.

22 e. Not more than 100 milligrams of opium per 100 milliliters or per 100 grams.

23 f. Not more than 0.5 milligram of difenoxin and not less than 25 micrograms of
24 atropine sulfate per dosage unit.

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

30 a. Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (also
31 referred to as BRV; UCB-34714; Briviact) (including its salts).

- 1 b. Cenobamate [(1R)-1-(2-chlorophenyl)-2-(tetrazol-2-yl)ethyl] carbamate; 2H-
2 tetrazole-2-ethanol, alpha-(2-chlorophenyl)-, carbamate (ester), (alphaR)-;
3 carbamic acid (R)-(+)-1-(2-chlorophenyl)-2-(2H-tetrazol-2-yl)ethyl ester).
4 c. Ezogabine N-[2-amino-4-(4-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester.
5 e.d. Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide].
6 d.e. Lasmiditan [2,4,6-trifluoro-N-(6-(1-methylpiperidine-4-carbonyl)pyridine-2-yl-
7 benzamide].
8 f. Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid].
9 e. ~~Approved cannabidiol drugs. A drug product in finished dosage formulation that~~
10 ~~has been approved by the federal food and drug administration, which contains~~
11 ~~cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-~~
12 ~~1,3-benzenediol) derived from cannabis and no more than 0.1 percent weight for~~
13 ~~weight residual tetrahydrocannabinols.~~
14 f.g. Gabapentin [2-[1-(aminomethyl) cyclohexyl] acetic acid].
15 6. Stimulants. Unless specifically exempted or excluded or unless listed in another
16 schedule, any material, compound, mixture, or preparation containing any quantity of
17 the following substances having a stimulant effect on the central nervous system,
18 including their salts, isomers, and salts of isomers: Pyrovalerone.

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