Sixty-seventh Legislative Assembly of North Dakota

SENATE BILL NO. 2059

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

6

7

8

9

10

11

12

13

14

15

16

17

18

19

20

21

22

BE IT ENACTED BY THE LEGISLATIVE ASSEMBLY OF NORTH DAKOTA:

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

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

from the seeds of the plant, fiber produced from the stalks, oil or cake made from the seeds of the plant, any other compound, manufacture, salt, derivative, mixture, or preparation of mature stalks (except the resin extracted therefrom),

Legislative Assembly 1 fiber, oil, or cake, or the sterilized seed of the plant which is incapable of 2 germination; 3 b. Hemp as defined in chapter 4.1-18.1; or 4 A prescription drug approved by the United States food and drug administration 5 under section 505 of the Federal Food, Drug, and Cosmetic Act [21 U.S.C. 355]. 6 SECTION 2. AMENDMENT. Section 19-03.1-05 of the North Dakota Century Code is 7 amended and reenacted as follows: 8 19-03.1-05. Schedule I. 9 1. The controlled substances listed in this section are included in schedule I. 10 2. Schedule I consists of the drugs and other substances, by whatever official name, 11 common or usual name, chemical name, or brand name designated, listed in this 12 section. 13 Opiates. Unless specifically excepted or unless listed in another schedule, any of the 3. 14 following opiates, including their isomers, esters, ethers, salts, and salts of isomers, 15 esters, and ethers, whenever the existence of those isomers, esters, ethers, and salts 16 is possible within the specific chemical designation: 17 Acetylmethadol. a. 18 b. Allylprodine. 19 Alphacetylmethadol. C. 20 d. Alphameprodine. 21 e. Alphamethadol. 22 f. Benzethidine. 23 Betacetylmethadol. g. 24 h. Betameprodine. 25 i. Betamethadol. 26 Betaprodine. j. 27 k. Brorphine. 28 Clonitazene. 29 Dextromoramide. ŀm. 30 m.n. Diampromide. 31 Diethylthiambutene. n.o.

Sixty-seventh

```
1
                        Difenoxin.
                <del>0.</del>p.
 2
                        Dimenoxadol.
                <del>p.</del>q.
 3
                <del>q.</del>r.
                        Dimepheptanol.
 4
                <del>r.</del>s.
                        Dimethylthiambutene.
 5
                <del>s.</del>t.
                        Dioxaphetyl butyrate.
 6
                <del>t.</del>u.
                        Dipipanone.
 7
                        Ethylmethylthiambutene.
                <del>U.</del>V.
 8
                        Etonitazene.
                <del>∀.</del><u>W.</u>
 9
               ₩.<u>X.</u>
                        Etoxeridine.
10
                <del>X.</del><u>y.</u>
                        Furethidine.
11
                <del>∀.</del>Z.
                        Hydroxypethidine.
12
              <del>z.</del>aa.
                        Isotonitazene.
13
              aa.bb. Ketobemidone.
14
              aa.bb.cc.
                                Levomoramide.
15
              bb.cc.dd.
                                Levophenacylmorphan.
16
                                Morpheridine.
              cc.dd.ee.
17
              dd.ee.ff.MPPP (also known as 1-methyl-4-phenyl-4-propionoxypiperidine).
18
              ee.ff.gg Noracymethadol.
19
              ff.gg.hh.Norlevorphanol.
20
              gg.hh.ii. Normethadone.
21
              hh.<u>ii.jj.</u> Norpipanone.
22
              ii.jj.kk. PEPAP (1-(2-Phenylethyl)-4-Phenyl-4-acetoxypiperidine).
23
              ij.kk.ll. Phenadoxone.
24
              kk.<u>ll.mm.</u>
                                Phenampromide.
25
              H.mm.nn.
                                Phenomorphan.
26
                                Phenoperidine.
              mm.nn.oo.
27
                                Piritramide.
              nn.<u>oo.</u>pp.
28
                                Proheptazine.
              <del>00.pp.qq.</del>
29
                                Properidine.
              pp.qq.rr.
30
              qq.<u>rr.ss.</u>Propiram.
31
              rr.ss.tt. Racemoramide.
```

1	ss.<u>tt.</u>uu.	ss. <u>tt.uu.</u> Tilidine.					
2	tt. <u>uu.</u> vv.	<u>vv.</u> Trimeperidine.					
3	uu.<u>vv.</u>wv	<u>v.</u>	3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also				
4	1	known	as U-47700).				
5	VV.<u>WW.</u>XX	<u>(.</u>	1-cyclohexyl-4-(1,2-diphenylethyl)piperazine (also know as MT-45).				
6	ww.xx.y	<u>/.</u>	3,4-dichloro- <i>N</i> -{[1-(dimethylamino)cyclohexyl]methyl}benzamide (also				
7]	known	as AH-7921).				
8	yy. zz.	Zipepr	ol.				
9	xx. <u>zz.</u> aa	<u>a.</u>	Fentanyl derivatives. Unless specifically excepted or unless listed in				
10	;	anothe	er schedule or are not FDA approved drugs, and are derived from N-(1-(2-				
11		Pheny	lethyl)-4-piperidinyl)-N-phenylpropanamide (Fentanyl) by any substitution				
12	(on or r	replacement of the phenethyl group, any substitution on the piperidine ring,				
13	;	any su	ibstitution on or replacement of the propanamide group, any substitution on				
14	f	the an	ilido phenyl group, or any combination of the above. Examples include:				
15	((1) N	I-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide (also known				
16		а	s Acetyl-alpha-methylfentanyl).				
17	((2) N	I-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide; 1-(1-methyl-				
18		2	-phenylethyl)-4-(N-propanilido)piperidine (also known as Alpha-				
19		n	nethylfentanyl).				
20	((3) N	I-[1-methyl-2-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (also				
21		k	nown as Alpha-methylthiofentanyl).				
22	((4) N	I-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-N-phenylpropanamide (also				
23		k	nown as Beta-hydroxyfentanyl).				
24	((5) N	I-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide				
25		(8	also known as Beta-hydroxy-3-methylfentanyl).				
26	((6) N	I-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (also				
27		k	nown as 3-Methylfentanyl).				
28	((7) N	I-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (also				
29		k	nown as 3-Methylthiofentanyl).				
30	((8) N	I-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl]propanamide (also				
31		k	nown as Para-fluorofentanyl).				

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

1		(24)	N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide (also known			
2				as Para-chloroisobutyryl Fentanyl).			
3		(25)	N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known			
4				as Para-methoxybutyryl Fentanyl).			
5		(26)	N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide (also known as			
6				Para-fluorobutyryl Fentanyl).			
7		(<u>27)</u>	N-(1-(2-fluorophenethyl)piperidin-4-yl)-N-(2-fluorophenyl)propionamide (also			
8				known as 2'-fluoro Ortho-fluorofentanyl).			
9		(<u>28)</u>	N-(2-methylphenyl)-N-(1-phenethylpiperidin-4-yl)acetamide (also known as			
10				Ortho-methyl Acetylfentanyl).			
11		(<u>29)</u>	N-(1-phenethylpiperidin-4-yl)-N,3-diphenylpropanamide (also known as			
12				Beta'-phenyl Fentanyl and Hydrocinnamoyl Fentanyl).			
13		(<u>30)</u>	N-(1-phenethylpiperidin-4-yl)-N-phenylthiophene-2-carboxamide (also			
14				known as Thiofuranyl Fentanyl).			
15		(<u>31)</u>	(E)-N-(1-phenethylpiperidin-4-yl)-N-phenylbut-2-enamide (also known as			
16				Crotonyl Fentanyl).			
17	4.	Opiu	ium derivatives. Unless specifically excepted or unless listed in another schedule,				
18		any	any of the following opium derivatives, its salts, isomers, and salts of isomers				
19		whenever the existence of such salts, isomers, and salts of isomers is possible within					
20		the specific chemical designation:					
21		a.	a. Acetorphine.				
22		b.	Acet	tyldihydrocodeine.			
23		C.	Ben	zylmorphine.			
24		d.	Cod	eine methylbromide.			
25		e.	Cod	eine-N-Oxide.			
26		f.	Сур	renorphine.			
27		g.	Des	omorphine.			
28		h.	Dihy	dromorphine.			
29		i.	Drot	ebanol.			
30		j.	Etor	phine (except hydrochloride salt).			
31		k.	k. Heroin.				

I.

Hydromorphinol.

1

2 Methyldesorphine. m. 3 n. Methyldihydromorphine. 4 Morphine methylbromide. 0. 5 Morphine methylsulfonate. p. 6 Morphine-N-Oxide. q. 7 Myrophine. r. 8 Nicocodeine. S. 9 t. Nicomorphine. 10 Normorphine. u. 11 ٧. Pholcodine. 12 Thebacon. W. 13 Hallucinogenic substances. Unless specifically excepted or unless listed in another 5. 14 schedule, any material, compound, mixture, or preparation containing any quantity of 15 the following hallucinogenic substances, including their salts, isomers, and salts of 16 isomers whenever the existence of those salts, isomers, and salts of isomers is 17 possible within the specific chemical designation (for purposes of this subsection only, 18 the term "isomer" includes the optical, position, and geometric isomers): 19 Alpha-ethyltryptamine, its optical isomers, salts, and salts of isomers (also known a. 20 as etryptamine; a-ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole). 21 b. Alpha-methyltryptamine. 22 4-methoxyamphetamine (also known as 4-methoxy-a-methylphenethylamine; C. 23 paramethoxyamphetamine; PMA). 24 d. N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-alpha-25 methyl-3,4(methylenedioxy)phenylamine, and N-hydroxy MDA. 26 Hashish. e. 27 f. Ibogaine (also known as 7-Ethyl-6, 6B, 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-28 6, 9-methano-5 H-pyrido [1', 2':1,2] azepino (5,4-b) indole; Tabernanthe iboga). 29 Lysergic acid diethylamide. g. 30 h. Marijuana.

1 Parahexyl (also known as 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro- 6,6,9-trimethyl-2 6H-dibenzol[b,d]pyran; Synhexyl). 3 j. Peyote (all parts of the plant presently classified botanically as Lophophora 4 williamsii Lemaire, whether growing or not, the seeds thereof, any extract from 5 any part of such plant, and every compound, manufacture, salts, derivative, 6 mixture, or preparation of such plant, its seeds, or its extracts). 7 N-ethyl-3-piperidyl benzilate. k. 8 Ι. N-methyl-3-piperidyl benzilate. 9 m. Psilocybin. 10 Tetrahydrocannabinols, meaning tetrahydrocannabinols naturally contained in a n. 11 plant of the genus Cannabis (cannabis plant), as well as synthetic equivalents of 12 the substances contained in the cannabis plant, or in the resinous extractives of 13 such plant, including synthetic substances, derivatives, and their isomers with 14 similar chemical structure and pharmacological activity to those substances 15 contained in the plant; excluding tetrahydrocannabinols found in hemp as defined 16 in title 4.1 chapter 4.1-18.1; such as the following: 17 (1) Delta-1 cis or trans tetrahydrocannabinol, and their optical isomers. Other 18 names: Delta-9-tetrahydrocannabinol. 19 (2) Delta-6 cis or trans tetrahydrocannabinol, and their optical isomers. 20 Delta-3,4 cis or trans tetrahydrocannabinol, and its optical isomers. (3) 21 (Since nomenclature of these substances is not internationally standardized, 22 compounds of these structures, regardless of numerical designation of atomic 23 positions covered.) 24 Cannabinoids, synthetic. It includes the chemicals and chemical groups listed Ο. 25 below, including their homologues, salts, isomers, and salts of isomers. The term 26 "isomer" includes the optical, position, and geometric isomers. 27 (1) Indole carboxaldehydes. Any compound structurally derived from 1H-indole-28 3-carboxaldehyde or 1H-2-carboxaldehyde substituted in both of the 29 following ways: at the nitrogen atom of the indole ring by an alkyl, haloalkyl, 30 cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-31 piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,

1	1-(N-met	hyl-3- morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo			
2	benzyl gr	benzyl group; and, at the hydrogen of the carboxaldehyde by a phenyl,			
3	benzyl, c	benzyl, cumyl, naphthyl, adamantyl, cyclopropyl, pyrrolidinyl, piperazinyl, or			
4	propiona	Idehyde group whether or not the compound is further modified to			
5	any exte	nt in the following ways:			
6	(a) Sul	bstitution to the indole ring to any extent; or			
7	(b) Sul	bstitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,			
8	сус	clopropyl, pyrrolidinyl, piperazinyl, or propionaldehyde group to any			
9	ext	ent; or			
10	(c) An	itrogen heterocyclic analog of the indole ring; or			
11	(d) An	itrogen heterocyclic analog of the phenyl, benzyl, naphthyl,			
12	ada	amantyl, or cyclopropyl ring.			
13	(e) Exa	amples include:			
14	[1]	1-Pentyl-3-(1-naphthoyl)indole - Other names: JWH-018 and			
15		AM-678.			
16	[2]	1-Butyl-3-(1-naphthoyl)indole - Other names: JWH-073.			
17	[3]	1-Pentyl-3-(4-methoxy-1-naphthoyl)indole - Other names:			
18		JWH-081.			
19	[4]	1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole - Other names:			
20		JWH-200.			
21	[5]	1-Propyl-2-methyl-3-(1-naphthoyl)indole - Other names:			
22		JWH-015.			
23	[6]	1-Hexyl-3-(1-naphthoyl)indole - Other names: JWH-019.			
24	[7]	1-Pentyl-3-(4-methyl-1-naphthoyl)indole - Other names:			
25		JWH-122.			
26	[8]	1-Pentyl-3-(4-ethyl-1-naphthoyl)indole - Other names: JWH-210.			
27	[9]	1-Pentyl-3-(4-chloro-1-naphthoyl)indole - Other names:			
28		JWH-398.			
29	[10]	1-(5-fluoropentyl)-3-(1-naphthoyl)indole - Other names:			
30		AM-2201.			

1		[11]	1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole - Other
2			names: RCS-8.
3		[12]	1-Pentyl-3-(2-methoxyphenylacetyl)indole - Other names:
4			JWH-250.
5		[13]	1-Pentyl-3-(2-methylphenylacetyl)indole - Other names:
6			JWH-251.
7		[14]	1-Pentyl-3-(2-chlorophenylacetyl)indole - Other names: JWH-
8			203.
9		[15]	1-Pentyl-3-(4-methoxybenzoyl)indole - Other names: RCS-4.
10		[16]	(1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole) - Other names:
11			AM-694.
12		[17]	(4-Methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-
13			yl]methanone - Other names: WIN 48,098 and Pravadoline.
14		[18]	(1-Pentylindol-3-yl)-(2,2,3,3-tetramethylcyclopropyl)methanone
15			Other names: UR-144.
16		[19]	(1-(5-fluoropentyl)indol-3-yl)-(2,2,3,3-
17			tetramethylcyclopropyl)methanone - Other names: XLR-11.
18		[20]	(1-(2-morpholin-4-ylethyl)-1H-indol-3-yl)-(2,2,3,3-
19			tetramethylcyclopropyl)methanone - Other names: A-796,260.
20		[21]	(1-(5-fluoropentyl)-1H-indazol-3-yl)(naphthalen-1-yl)methanone
21			Other names: THJ-2201.
22		[22]	1-naphthalenyl(1-pentyl-1H-indazol-3-yl)-methanone Other
23			names: THJ-018.
24		[23]	(1-(5-fluoropentyl)-1H-benzo[d]imidazol-2-yl)(naphthalen-1-
25			yl)methanone - Other names: FUBIMINA.
26		[24]	1-[(N-methylpiperidin-2-yl)methyl]-3-(adamant-1-oyl) indole -
27			Other names: AM-1248.
28		[25]	1-Pentyl-3-(1-adamantoyl)indole - Other names: AB-001 and
29			JWH-018 adamantyl analog.
30	(2)	Indole car	boxamides. Any compound structurally derived from 1H-indole-3-
31		carboxam	ide or 1H-2-carboxamide substituted in both of the following ways:

1	at th	e nitro	ogen atom of the indole ring by an alkyl, haloalkyl, cyanoalkyl,		
2	alkei	nyl, c	ycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-piperidinyl)methyl,		
3	2-(4-	2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-			
4	morp	holin	yl)methyl, tetrahydropyranylmethyl, benzyl, or halo benzyl group;		
5	and,	at the	e nitrogen of the carboxamide by a phenyl, benzyl, cumyl,		
6	naph	ıthyl,	adamantyl, cyclopropyl, or propionaldehyde group whether or not		
7	the c	ompo	ound is further modified to any extent in the following ways:		
8	(a)	Sub	stitution to the indole ring to any extent; or		
9	(b)	Sub	stitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,		
10		cycl	opropyl, or propionaldehyde group to any extent; or		
11	(c)	A ni	trogen heterocyclic analog of the indole ring; or		
12	(d)	A ni	trogen heterocyclic analog of the phenyl, benzyl, naphthyl,		
13		ada	mantyl, or cyclopropyl ring.		
14	(e)	Exa	mples include:		
15		[1]	N-Adamantyl-1-pentyl-1H-indole-3-carboxamide - Other names:		
16			JWH-018 adamantyl carboxamide, APICA, SDB-001, and 2NE1.		
17		[2]	N-Adamantyl-1-fluoropentylindole-3-carboxamide - Other names:		
18			STS-135.		
19		[3]	N-Adamantyl-1-pentyl-1H-Indazole-3-carboxamide - Other		
20			names: AKB 48 and APINACA.		
21		[4]	N-1-naphthalenyl-1-pentyl-1H-indole-3-carboxamide - Other		
22			names: NNEI and MN-24.		
23		[5]	N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-		
24			carboxamide - Other names: ADBICA.		
25		[6]	(S)-N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-		
26			3-carboxamide - Other names: AB-PINACA.		
27		[7]	N-[(1S)-1-(aminocarbonyl)-2-methylpropyl]-1-[(4-		
28			fluorophenyl)methyl]-1H-indazole-3-carboxamide - Other names:		
29			AB-FUBINACA.		

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

1		[19]	Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-
2			dimethylbutanoate - Other names: MDMB-CHMICA and
3			MMB-CHMINACA.
4		[20]	Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-
5			dimethylbutanoate - Other names: MDMB-FUBINACA.
6		[21]	1-(4-cyanobutyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carbox
7			amide - Other names: 4-CN-CUMYL-BUTINACA; 4-cyano-
8			CUMYL-BUTINACA; 4-CN-CUMYL BINACA; CUMYL-4CN -
9			BINACA; SGT-78.
10		[22]	methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-
11			3-methylbutanoate - Other names: MMB-CHMICA, AMB-
12			CHMICA.
13		[23]	1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-pyrrolo[2,3-b]pyrid
14			ine-3-carboxamide - Other names: 5F-CUMYL-P7AICA.
15		[24]	ethyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-
16			dimethylbutanoate - Other names: 5F-EDMB-PINACA.
17		[25]	methyl 2-(1-(5-fluoropentyl)-1H-indole-3-carboxamido)-3,3-
18			dimethylbutanoate - Other names: 5F-MDMB-PICA.
19		[26]	1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-
20			carboxamide - Other names: 5F-CUMYL-PINACA, SGT-25).
21		[27]	(1-(4-fluorobenzyl)-1H-indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)
22			methanone - Other names: FUB-144.
23	(3)	Indole car	boxylic acids. Any compound structurally derived from 1H-indole-
24		3-carboxy	lic acid or 1H-2-carboxylic acid substituted in both of the following
25		ways: at th	ne nitrogen atom of the indole ring by an alkyl, haloalkyl,
26		cyanoalky	l, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-2-
27		piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-pyrrolidinyl)methyl,
28		1-(N-meth	yl-3- morpholinyl)methyl, tetrahydropyranylmethyl, benzyl, or halo
29		benzyl gro	oup; and, at the hydroxyl group of the carboxylic acid by a phenyl,
30		benzyl, cu	myl, naphthyl, adamantyl, cyclopropyl, or propionaldehyde group

1		whet	her o	r not the compound is further modified to any extent in the
2		follo	wing v	ways:
3		(a)	Sub	stitution to the indole ring to any extent; or
4		(b)	Sub	stitution to the phenyl, benzyl, cumyl, naphthyl, adamantyl,
5			cycl	opropyl, propionaldehyde group to any extent; or
6		(c)	A nit	trogen heterocyclic analog of the indole ring; or
7		(d)	A nit	trogen heterocyclic analog of the phenyl, benzyl, naphthyl,
8			adaı	mantyl, or cyclopropyl ring.
9		(e)	Exa	mples include:
10			[1]	1-(cyclohexylmethyl)-1H-indole-3-carboxylic acid 8-quinolinyl
11				ester - Other names: BB-22 and QUCHIC.
12			[2]	naphthalen-1-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate -
13				Other names: FDU-PB-22.
14			[3]	1-pentyl-1H-indole-3-carboxylic acid 8-quinolinyl ester - Other
15				names: PB-22 and QUPIC.
16			[4]	1-(5-Fluoropentyl)-1H-indole-3-carboxylic acid 8-quinolinyl ester
17				Other names: 5-Fluoro PB-22 and 5F-PB-22.
18			[5]	quinolin-8-yl-1-(4-fluorobenzyl)-1H-indole-3-carboxylate - Other
19				names: FUB-PB-22.
20			[6]	naphthalen-1-yl 1-(5-fluoropentyl)-1H-indole-3-carboxylate -
21				Other names: NM2201 and CBL2201.
22	(4)	Napl	nthyln	nethylindoles. Any compound containing a 1H-indol-3-yl-(1-
23		naph	thyl)r	methane structure with substitution at the nitrogen atom of the
24		indol	e ring	by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
25		cyclo	alkyle	ethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-
26		(N-m	ethyl-	-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or
27		(tetra	ahydro	opyran-4-yl)methyl group whether or not further substituted in the
28		indol	e ring	g to any extent and whether or not substituted in the naphthyl ring
29		to ar	y ext	ent. Examples include:
30		(a)	1-Pe	entyl-1H-indol-3-yl-(1-naphthyl)methane - Other names: JWH-175.

1		(b)	1-Pentyl-1H-indol-3-yl-(4-methyl-1-naphthyl)methane - Other names:
2			JWH-184.
3	(5)	Napl	nthoylpyrroles. Any compound containing a 3-(1-naphthoyl)pyrrole
4		struc	cture with substitution at the nitrogen atom of the pyrrole ring by an
5		alkyl	, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-
6		meth	nyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-(N-methyl-2-
7		pyrro	olidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-
8		yl)me	ethyl group whether or not further substituted in the pyrrole ring to any
9		exte	nt, whether or not substituted in the naphthyl ring to any extent.
0		Exar	mples include: (5-(2-fluorophenyl)-1-pentylpyrrol-3-yl)-naphthalen-1-
11		ylme	thanone - Other names: JWH-307.
2	(6)	Napl	nthylmethylindenes. Any compound containing a naphthylideneindene
3		struc	cture with substitution at the 3-position of the indene ring by an alkyl,
4		halo	alkyl, cyanoalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl, 1-(N-methyl-
5		2-pip	peridinyl)methyl, 2 (4 morpholinyl)ethyl, 1-(N-methyl-2-
16		pyrro	olidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or (tetrahydropyran-4-
7		yl)m	ethyl group whether or not further substituted in the indene ring to any
8		exte	nt, whether or not substituted in the naphthyl ring to any extent.
9		Exar	mples include: E-1-[1-(1-Naphthalenylmethylene)-1H-inden-3-yl]pentane
20		- Oth	ner names: JWH-176.
21	(7)	Cycl	ohexylphenols. Any compound containing a 2-(3-
22		hydr	oxycyclohexyl)phenol structure with substitution at the 5-position of the
23		pher	nolic ring by an alkyl, haloalkyl, cyanoalkyl, alkenyl, cycloalkylmethyl,
24		cyclo	palkylethyl, 1-(N-methyl-2-piperidinyl)methyl, 2-(4-morpholinyl)ethyl, 1-
25		(N-m	nethyl-2-pyrrolidinyl)methyl, 1-(N-methyl-3-morpholinyl)methyl, or
26		(tetra	ahydropyran-4-yl)methyl group whether or not substituted in the
27		cyclo	phexyl ring to any extent. Examples include:
28		(a)	5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
29			names: CP 47,497.
30		(b)	5-(1,1-dimethyloctyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol - Other
31			names: Cannabicyclohexanol and CP 47,497 C8 homologue.

1			(c)	5-(1,1-dimethylheptyl)-2-[(1R,2R)-5-hydroxy-2-(3-		
2				hydroxypropyl)cyclohexyl]-phenol - Other names: CP 55,940.		
3		(8)	Othe	rs specifically named:		
4			(a)	(6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-		
5				6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names: HU-210.		
6			(b)	(6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-		
7				6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol - Other names:		
8				Dexanabinol and HU-211.		
9			(c)	2,3-Dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3-de]-1,4-		
10				benzoxazin-6-yl]-1-napthalenylmethanone - Other names:		
11				WIN 55,212-2.		
12			(d)	Naphthalen-1-yl-(4-pentyloxynaphthalen-1-yl)methanone - Other		
13				names: CB-13.		
14	p.	Sub	stitute	d phenethylamines. This includes any compound, unless specifically		
15		exce	excepted, specifically named in this schedule, or listed under a different			
16		sche	schedule, structurally derived from phenylethan-2-amine by substitution on the			
17		phei	phenyl ring in any of the following ways, that is to say, by substitution with a fused			
18		metl	methylenedioxy ring, fused furan ring, or fused tetrahydrofuran ring; by			
19		subs	substitution with two alkoxy groups; by substitution with one alkoxy and either			
20		one	fused	furan, tetrahydrofuran, or tetrahydropyran ring system; or by		
21		subs	stitutio	n with two fused ring systems from any combination of the furan,		
22		tetra	hydro	furan, or tetrahydropyran ring systems.		
23		(1)	Whe	ther or not the compound is further modified in any of the following		
24			ways	s, that is to say:		
25			(a)	By substitution of phenyl ring by any halo, hydroxyl, alkyl,		
26				trifluoromethyl, alkoxy, or alkylthio groups;		
27			(b)	By substitution at the 2-position by any alkyl groups; or		
28			(c)	By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl,		
29				hydroxybenzyl, methylenedioxybenzyl, or methoxybenzyl groups.		
30		(2)	Exan	nples include:		

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

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

1		(dd)	2,5-dimethoxy-4-ethylamphetamine (also known as DOET).
2		(ee)	2,5-dimethoxy-4-(n)-propylthiophenethylamine (also known as 2C-T-
3			7).
4		(ff)	5-methoxy-3,4-methylenedioxy-amphetamine.
5		(gg)	4-methyl-2,5-dimethoxy-amphetamine (also known as 4-methyl-2,5-
6			dimethoxy-a-methylphenethylamine; DOM and STP).
7		(hh)	3,4-methylenedioxy amphetamine (also known as MDA).
8		(ii)	3,4-methylenedioxymethamphetamine (also known as MDMA).
9		(jj)	3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-
0			alpha-methyl-3,4(methylenedioxy)phenethylamine, MDE, MDEA).
11		(kk)	3,4,5-trimethoxy amphetamine.
2		(II)	Mescaline (also known as 3,4,5-trimethoxyphenethylamine).
3		(mm)	1-(4-methoxyphenyl)-N-methylpropan-2-amine (also known as para-
4			methoxymethamphetamine and PMMA).
5	q.	Substitu	ted tryptamines. This includes any compound, unless specifically
6		excepted	d, specifically named in this schedule, or listed under a different
7		schedule	e, structurally derived from 2-(1H-indol-3-yl)ethanamine (i.e., tryptamine)
8		by mond	o- or di-substitution of the amine nitrogen with alkyl or alkenyl groups or
9		by inclus	sion of the amino nitrogen atom in a cyclic structure whether or not the
20		compou	nd is further substituted at the alpha-position with an alkyl group or
21		whether	or not further substituted on the indole ring to any extent with any alkyl,
22		alkoxy, h	nalo, hydroxyl, or acetoxy groups. Examples include:
23		(1) 5-n	nethoxy-N,N-diallyltryptamine (also known as 5-MeO-DALT).
24		(2) 4-a	cetoxy-N,N-dimethyltryptamine (also known as 4-AcO-DMT or O-
25		Ace	etylpsilocin).
26		(3) 4-h	ydroxy-N-methyl-N-ethyltryptamine (also known as 4-HO-MET).
27		(4) 4-h	ydroxy-N,N-diisopropyltryptamine (also known as 4-HO-DIPT).
28		(5) 5-n	nethoxy-N-methyl-N-isopropyltryptamine (also known as 5-MeO-MiPT).
<u> 2</u> 9		(6) 5-n	nethoxy-N,N-dimethyltryptamine (also known as 5-MeO-DMT).

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

1		<u>e.</u>	<u>Etizo</u>	olam.	
2		<u>f.</u>	<u>Flua</u>	lprazo	olam.
3		<u>g.</u>	<u>Flub</u>	romaz	<u>repam.</u>
4		<u>h.</u>	Flub	romaz	zolam.
5		<u>i.</u>	<u>Adin</u>	azola	<u>m.</u>
6	7.	Stin	nulant	s. Unl	ess specifically excepted or unless listed in another schedule, any
7		mat	erial,	compo	ound, mixture, or preparation which contains any quantity of the
8		follo	wing	substa	ances having a stimulant effect on the central nervous system,
9		incl	uding	its sal	ts, isomers, and salts of isomers:
10		a.	Amii	norex	(also known as 2-amino-5-phenyl-2-oxazoline, or 4,5-dihydro-5-phenyl-
11			2-ox	azolaı	mine).
12		b.	Cath	ninone	
13		C.	Sub	stitute	d cathinones. Any compound, material, mixture, preparation, or other
14			prod	luct, u	nless listed in another schedule or an approved food and drug
15			adm	inistra	tion drug (e.g., buproprion, pyrovalerone), structurally derived from 2-
16			amir	noprop	pan-1-one by substitution at the 1-position with either phenyl, naphthyl,
17			or th	iophe	ne ring systems, whether or not the compound is further modified in
18			any	of the	following ways:
19			(1)	By su	ubstitution in the ring system to any extent with alkyl, alkylenedioxy,
20				alkox	y, haloalkyl, hydroxyl, or halide substituents, whether or not further
21				subs	tituted in the ring system by one or more other univalent substitutents;
22			(2)	By su	ubstitution at the 3-position with an acyclic alkyl substituent;
23			(3)	By su	ubstitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl, or
24				meth	oxybenzyl groups; or
25			(4)	By in	clusion of the 2-amino nitrogen atom in a cyclic structure.
26				Some	e trade or other names:
27				(a)	3,4-Methylenedioxy-alpha-pyrrolidinopropiophenone (also known as
28					MDPPP).
29				(b)	3,4-Methylenedioxy-N-ethylcathinone (also known as Ethylone,
30					MDEC, or bk-MDEA).

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

1			<u>(aa)</u>	N-ethylhexedrone.	
2			<u>(bb)</u>	alpha-pyrrolidinohexanophenone (also known as alpha-PHP).	
3			<u>(cc)</u>	4-methyl-alpha-ethylaminopentiophenone (also known as 4-MEAP).	
4			<u>(dd)</u>	4'-methyl-alpha-pyrrolidinohexiophenone (also known as MPHP).	
5			<u>(ee)</u>	alpha-pyrrolidinoheptaphenone (also known as PV8).	
6			<u>(ff)</u>	4-chloro-alpha-pyrrolidinovalerophenone (also known 4-chloro-alpha-	
7				PVP).	
8		d.	Fenethylli	ne.	
9		e.	Fluoroam	phetamine.	
0		f.	Fluoromet	thamphetamine.	
11		g.	(±)cis-4-m	nethylaminorex (also known as (±)cis-4,5-dihydro-4-methyl-5-phenyl-2-	
2			oxazolam	ine).	
3		h.	N-Benzylp	piperazine (also known as BZP, 1-benzylpiperazine).	
4		i.	N-ethylam	phetamine.	
5		j.	N, N-dime	ethylamphetamine (also known as N,N-alpha-trimethyl-	
6			benzenee	thanamine; 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	9 19-03.1-07. Schedule II.				
20	1.	The controlled substances listed in this section are included in schedule II.			
21	2.	Schedule II consists of the drugs and other substances, by whatever official name,			
22		common or usual name, chemical name, or brand name designated, listed in this			
23		sec	tion.		
24	3.	Sub	stances, ve	egetable origin or chemical synthesis. Unless specifically excepted or	
25		unle	ess listed in	another schedule, any of the following substances whether produced	
26		directly or indirectly by extraction from substances of vegetable origin, or			
27		independently by means of chemical synthesis, or by a combination of extraction and			
28		che	mical synth	esis:	
29		a.	Opium an	d opiate, and any salt, compound, derivative, or preparation of opium or	
30	I		opiate, ex	cluding apomorphine, thebaine-derived butorphanol, dextrorphan,	
31			nalbuphin	e, <u>naldemedine,</u> nalmefene, <u>naloxegol,</u> naloxone, <u>6 beta-naltrexol,</u> and	

1 naltrexone, and samidorphan and their respective salts, but including the 2 following: 3 (1) Codeine. 4 (2) Dihydroetorphine. 5 (3) Ethylmorphine. 6 (4) Etorphine hydrochloride. 7 (5)Granulated opium. 8 (6) Hydrocodone. 9 (7) Hydromorphone. 10 (8) Metopon. 11 (9)Morphine. 12 (10)Noroxymorphone. 13 (11)Opium extracts. 14 (11)(12) Opium fluid. 15 (12)(13) Oripavine. 16 (13)(14) Oxycodone. 17 (14)(15) Oxymorphone. 18 (15)(16) Powder opium. 19 (16)(17) Raw opium. 20 (17)(18) Thebaine. 21 (18)(19) Tincture of opium. 22 b. Any salt, compound, derivative, or preparation thereof which is chemically 23 equivalent or identical with any of the substances referred to in subdivision a, but 24 not including the isoquinoline alkaloids of opium. 25 Opium poppy and poppy straw. C. 26 Coca leaves and any salt, compound, derivative, or preparation of coca leaves, d. 27 including cocaine and ecgonine and their salts, isomers, derivatives, and salts of 28 isomers and derivatives, and any salt, compound, derivative, or preparation 29 thereof that is chemically equivalent or identical with any of these substances. 30 except that the nondosage substances must include decocainized coca leaves or 31 extractions of coca leaves which do not contain cocaine or ecgonine.

1 Concentrate of poppy straw (the crude extract of poppy straw in either liquid, 2 solid, or powder form which contains the phenanthrine alkaloids of the opium 3 poppy). 4 4. Opiates. Unless specifically excepted or unless 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, dextrophan and 8 levopropoxyphene excepted: 9 Alfentanil. a. 10 Alphaprodine. b. 11 C. Anileridine. 12 d. Bezitramide. 13 Bulk dextropropoxyphene (nondosage forms). e. 14 f. Carfentanil. 15 g. Dihydrocodeine. 16 h. Diphenoxylate. 17 i. Fentanyl. 18 j. Isomethadone. 19 k. Levo-alphaacetylmethadol (LAAM). 20 Ι. Levomethorphan. 21 m. Levorphanol. 22 Metazocine. n. 23 Methadone. 0. 24 p. Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane. 25 Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic q. 26 acid. 27 Oliceridine (N-[(3-methoxythiophen-2-yl)methyl] ({2-[(9R)-9-(pyridin-2-yl)-6r. 28 oxaspiro [4.5]decan-9-yl]ethyl})amine fumarate). 29 Pethidine (also known as meperidine). S. 30 s.t. Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine. 31 Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate. t.u.

- 1 <u>u.v.</u> Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid.
- 2 v.w. Phenazocine.
- $\frac{3}{\text{w.x.}}$ Priminodine.
- 4 x.y. Racemethorphan.
- 5 <u>y.z.</u> Racemorphan.
- 6 z.aa. Remifentanil.
- 7 aa.bb. Sufentanil.
- 8 <u>bb.cc.</u> Tapentadol.
- 9 <u>cc.dd.</u> Thiafentanil.

13

15

- 5. Stimulants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a stimulant effect on the central nervous system:
 - a. Amphetamine, its salts, optical isomers, and salts of its optical isomers.
- b. Lisdexamfetamine, its salts, isomers, and salts of isomers.
 - c. Methamphetamine, its salts, isomers, and salts of isomers.
- 16 d. Phenmetrazine and its salts.
- 17 e. Methylphenidate.
- 6. Depressants. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation which contains any quantity of the following substances having a depressant effect on the central nervous system, including its salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation:
- a. Amobarbital.
- b. Glutethimide.
- c. Pentobarbital.
- d. Phencyclidine.
- e. Secobarbital.
- 28 7. Hallucinogenic substances.
- 29 a. Nabilone [another name for nabilone (±)-trans-3-(1, 1-dimethylheptyl)-6, 6a, 7, 8, 10, 10a-hexahydro-1-hydroxy-6, 6-dimethyl-9Hdibenzo [b, d] pyran-9-one].

1		b.	Dronabinol [(-)-delta-9-trans tetrahydrocannabinol] in an oral solution in a drug					
2			product approved for marketing by the federal food and drug administration.					
3	8.	lmn	Immediate precursors. Unless specifically excepted or unless listed in another					
4		sch	edule, any material, compound, mixture, or preparation that contains any quantity					
5		of t	he following substances:					
6		a.	Immediate precursor to amphetamine and methamphetamine: Phenylacetone.					
7			Some trade or other names: phenyl-2-propanone; P2P, benzyl methyl ketone;					
8			methyl benzyl ketone.					
9		b.	Immediate precursors to phencyclidine (PCP):					
10			(1) 1-phenylcyclohexylamine.					
11	ı		(2) 1-piperidinocyclohexanecarbonitrile (PCC).					
12		C.	Immediate precursors to fentanyl: 4-anilino-N-phenethyl-4-piperidine (ANPP)					
13			(1) 4-anilino-N-phenethylpiperidine (ANPP).					
14			(2) N-phenyl-N-(piperidin-4-yl)propionamide (norfentanyl).					
15	SEC	SECTION 4. AMENDMENT. Section 19-03.1-11 of the North Dakota Century Code is						
16	amende	amended and reenacted as follows:						
17	19-0	9-03.1-11. Schedule IV.						
18	1.	The	e controlled substances listed in this section are included in schedule IV.					
19	2.	Sch	Schedule IV consists of the drugs and other substances, by whatever official name,					
20		con	nmon or usual name, chemical name, or brand name designated, listed in this					
21		sec	tion.					
22	3.	Nar	cotic drugs. Unless specifically excepted or unless listed in another schedule, any					
23		material, compound, mixture, or preparation containing any of the following narcotic						
24		drugs or their salts calculated as the free anhydrous base or alkaloid, in limited						
25		quantities as set forth below:						
26		a.	Not more than 1 milligram of difenoxin and not less than 25 micrograms of					
27			atropine sulfate per dosage unit.					
28		b.	Dextropropoxyphene (also known as alpha-(+)-4-dimethylamino- 1,2-diphenyl-3-					
29			methyl-2-propionoxybutane).					
30		C.	2-[(dimethylamino)methyl]-1-(3-methoxyphenyl)cyclohexanol, its salts, optical					
31			and geometric isomers and salts of these isomers including Tramadol.					

31

Fospropofol.

∀.Z.

1 Depressants. Unless specifically excepted or unless listed in another schedule, any 2 material, compound, mixture, or preparation containing any quantity of the following 3 substances, including their salts, isomers, and salts of isomers whenever the 4 existence of those salts, isomers, and salts of isomers is possible within the specific 5 chemical designation: 6 a. Alprazolam. 7 b. Alfaxalone. 8 Barbital. C. 9 d. Brexanolone. 10 <u>e.</u> Bromazepam. 11 e.f. Camazepam. 12 <u>f.g.</u> Carisoprodol. 13 g.h. Chloral betaine. 14 h.i. Chloral hydrate. 15 i.j. Chlordiazepoxide. 16 Clobazam. j.k. 17 k.l. Clonazepam. 18 l.m. Clorazepate. 19 Clotiazepam. m.n. 20 Cloxazolam. n.o. 21 о.р. Delorazepam. 22 Diazepam. p.g. 23 Dichloralphenazone. q.r. 24 r.s. Estazolam. 25 s.t. Ethchlorvynol. 26 Ethinamate. t.u. 27 U.<u>V.</u> Ethyl loflazepate. 28 Fludiazepam. ₩. 29 Flunitrazepam. ₩.<u>X.</u> 30 X.<u>y.</u> Flurazepam.

1 Halazepam. z.aa. 2 aa.bb. Haloxazolam. 3 bb.cc. Indiplon. 4 ec.dd. Ketazolam. 5 dd.ee. Lemborexant. 6 ff. Loprazolam. 7 ee.gg. Lorazepam. 8 ff.hh. Lorcaserin. 9 gg.ii. Lormetazepam. 10 hh.jj. Mebutamate. 11 ii.kk. Medazepam. 12]].[]. Meprobamate. 13 kk.mm. Methohexital. 14 Methylphenobarbital (also known as mephobarbital). <u>₩.nn.</u> 15 mm.oo. Midazolam. 16 nn.pp. Nimetazepam. 17 oo.qq. Nitrazepam. 18 pp.rr. Nordiazepam. 19 Oxazepam. qq.ss. 20 rr.tt. Oxazolam. 21 ss.uu. Paraldehyde. 22 tt.vv. Petrichloral. 23 uu.ww. Phenobarbital. 24 ₩.<u>XX.</u> Pinazepam. 25 ww.yy. Propofol. 26 Prazepam. XX.<u>ZZ.</u> 27 yy.aaa. Quazepam. 28 zz.bbb. Remimazolam. 29 Suvorexant. CCC. 30 aaa.ddd. Temazepam. 31 bbb.eee. Tetrazepam.

- 1 <u>ccc.fff.</u> Triazolam.
- 2 ddd.ggg. Zaleplon.
- 3 <u>eee.hhh.</u> Zolpidem.
- 4 fff.iii. Zopiclone.
- 5 5. Fenfluramine. Any material, compound, mixture, or preparation which contains any
- 6 quantity of the following substances, including its salts, isomers (whether optical,
- position, or geometric), and salts of such isomers, whenever the existence of such
- 8 salts, isomers, and salts of isomers is possible: Fenfluramine.
- 9 6. Stimulants. Unless specifically excepted or unless listed in another schedule, any
- material, compound, mixture, or preparation which contains any quantity of the
- following substances having a stimulant effect on the central nervous system,
- including its salts, isomers, and salts of isomers:
- 13 a. Cathine.
- b. Diethylpropion.
- 15 c. Fencamfamin.
- d. Fenproporex.
- 17 e. Mazindol.
- 18 f. Mefenorex.
- 19 g. Modafinil.
- 20 h. Pemoline (including organometallic complexes and chelates thereof).
- i. Phentermine.
- j. Pipradrol.
- k. Sibutramine.
- 24 I. <u>Solriamfetol.</u>
- 25 <u>m.</u> SPA ((-)-1-dimethylamino-1, 2-diphenylethane).
- 7. Other substances. Unless specifically excepted or unless listed in another schedule,
- any material, compound, mixture, or preparation which contains any quantity of:
- 28 a. Pentazocine, including its salts.
- b. Butorphanol, including its optical isomers.
- 30 c. Eluxadoline (5-[[[(2S)-2-amino-3-[4-aminocarbonyl)-2,6-dimethylphenyl]-1-
- 31 oxopropyl][(1S)-1-(4-phenyl-1H-imidazol-2-yl)ethyl]amino]methyl]-2-

3

4

5

6

7

8

9

12

13

17

18

19

20

21

22

23

24

25

26

27

28

29

30

- methoxybenzoic acid) (including its optical isomers) and its salts, isomers, and salts of isomers.
 - 8. The board may except by rule any compound, mixture, or preparation containing any depressant substance listed in subsection 2 from the application of all or any part of this chapter if the compound, mixture, or preparation contains one or more active medicinal ingredients not having a depressant effect on the central nervous system, and if the admixtures are included therein in combinations, quantity, proportion, or concentration that vitiate the potential for abuse of the substances which have a depressant effect on the central nervous system.

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

19-03.1-13. Schedule V.

- 1. The controlled substances listed in this section are included in schedule V.
- Schedule V consists of the drugs and other substances, by whatever official name,
 common or usual name, chemical name, or brand name designated, listed in this
 section.
 - Narcotic drugs. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation containing any of the following narcotic drugs and their salts.
 - 4. Narcotic drugs containing non-narcotic active medicinal ingredients. Any compound, mixture, or preparation containing any of the following narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid, in limited quantities as set forth below, which includes one or more non-narcotic active medicinal ingredients in sufficient proportion to confer upon the compound, mixture, or preparation valuable medicinal qualities other than those possessed by narcotic drugs alone.
 - a. Not more than 200 milligrams of codeine per 100 milliliters or per 100 grams.
 - Not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100 grams.
 - Not more than 100 milligrams of ethylmorphine per 100 milliliters or per
 100 grams.

31

1 Not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms 2 of atropine sulfate per dosage unit. 3 e. Not more than 100 milligrams of opium per 100 milliliters or per 100 grams. 4 f. Not more than 0.5 milligram of difenoxin and not less than 25 micrograms of 5 atropine sulfate per dosage unit. 6 5. Depressants. Unless specifically exempted or excluded or unless listed in another 7 schedule, any material, compound, mixture, or preparation that contains any quantity 8 of the following substances having a depressant effect on the central nervous system, 9 including its salts, isomers, and salts of isomers whenever the existence of such salts, 10 isomers, and salts of isomers is possible: 11 Brivaracetam ((2S)-2-[(4R)-2-oxo-4-propylpyrrolidin-1-yl]butanamide) (also 12 referred to as BRV; UCB-34714; Briviact) (including its salts). 13 b. Cenobamate [(1R)-1-(2-chlorophenyl)-2-(tetrazol-2-yl)ethyl] carbamate; 2H-14 tetrazole-2-ethanol, alpha-(2-chlorophenyl)-, carbamate (ester), (alphaR)-; 15 carbamic acid (R)-(+)-1-(2-chlorophenyl)-2-(2H-tetrazol-2-yl)ethyl ester). 16 Ezogabine N-[2-amino-4-(4-fluorobenzylamino)-phenyl]-carbamic acid ethyl ester. <u>C.</u> 17 c.d. Lacosamide [(R)-2-acetoamido-N-benzyl-3-methoxy-propionamide]. 18 <u>d.e.</u> Lasmiditan [2,4,6-trifluoro-N-(6-(1-methylpiperidine-4-carbonyl)pyridine-2-yl-19 benzamide]. 20 Pregabalin [(S)-3-(aminomethyl)-5-methylhexanoic acid]. <u>f.</u> 21 e. Approved cannabidiol drugs. A drug product in finished dosage formulation that 22 has been approved by the federal food and drug administration, which contains 23 cannabidiol (2-[1R-3-methyl-6R-(1-methylethenyl)-2-cyclohexen-1-yl]-5-pentyl-24 1,3-benzenediol) derived from cannabis and no more than 0.1 percent weight for 25 weight residual tetrahydrocannabinols. 26 Gabapentin [2-[1-(aminomethyl) cyclohexyl] acetic acid]. <u>f.g.</u> 27 6. Stimulants. Unless specifically exempted or excluded or unless listed in another 28 schedule, any material, compound, mixture, or preparation containing any quantity of 29 the following substances having a stimulant effect on the central nervous system. 30 including their salts, isomers, and salts of isomers: Pyrovalerone.

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