By: Miles

S.B. No. 1646

	A BILL TO BE ENTITLED
1	AN ACT
2	relating to the controlled substances listed in Penalty Groups 1,
3	2, and 2-A under the Texas Controlled Substances Act.
4	BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF TEXAS:
5	SECTION 1. Section 481.102, Health and Safety Code, is
6	amended to read as follows:
7	Sec. 481.102. PENALTY GROUP 1. Penalty Group 1 consists
8	of:
9	(1) the following opiates, including their isomers,
10	esters, ethers, salts, and salts of isomers, esters, and ethers,
11	unless specifically excepted, if the existence of these isomers,
12	esters, ethers, and salts is possible within the specific chemical
13	designation:
14	Alfentanil;
15	Allylprodine;
16	Alphacetylmethadol;
17	Benzethidine;
18	Betaprodine;
19	Clonitazene;
20	Diampromide;
21	Diethylthiambutene;
22	Difenoxin not listed in Penalty Group 3 or 4;
23	Dimenoxadol;
24	Dimethylthiambutene;

1	Dioxaphetyl butyrate;
2	Dipipanone;
3	Ethylmethylthiambutene;
4	Etonitazene;
5	Etoxeridine;
6	Furethidine;
7	Hydroxypethidine;
8	Ketobemidone;
9	Levophenacylmorphan;
10	Meprodine;
11	Methadol;
12	Moramide;
13	Morpheridine;
14	Noracymethadol;
15	Norlevorphanol;
16	Normethadone;
17	Norpipanone;
18	Phenadoxone;
19	Phenampromide;
20	Phenomorphan;
21	Phenoperidine;
22	Piritramide;
23	Proheptazine;
24	Properidine;
25	Propiram;
26	Sufentanil;
27	Tilidine; and

1	Trimeperidine;
2	(2) the following opium derivatives, their salts,
3	isomers, and salts of isomers, unless specifically excepted, if the
4	existence of these salts, isomers, and salts of isomers is possible
5	within the specific chemical designation:
6	Acetorphine;
7	Acetyldihydrocodeine;
8	Benzylmorphine;
9	Codeine methylbromide;
10	Codeine-N-Oxide;
11	Cyprenorphine;
12	Desomorphine;
13	Dihydromorphine;
14	Drotebanol;
15	Etorphine, except hydrochloride salt;
16	Heroin;
17	Hydromorphinol;
18	Methyldesorphine;
19	Methyldihydromorphine;
20	Monoacetylmorphine;
21	Morphine methylbromide;
22	Morphine methylsulfonate;
23	Morphine-N-Oxide;
24	Myrophine;
25	Nicocodeine;
26	Nicomorphine;
27	Normorphine;

1	Pholcodine; and
2	Thebacon;
3	(3) the following substances, however produced,
4	except those narcotic drugs listed in another group:
5	(A) Opium and opiate not listed in Penalty Group
6	3 or 4, and a salt, compound, derivative, or preparation of opium or
7	opiate, other than thebaine derived butorphanol, nalmefene and its
8	salts, naloxone and its salts, and naltrexone and its salts, but
9	including:
10	Codeine not listed in Penalty Group 3 or 4;
11	Dihydroetorphine;
12	Ethylmorphine not listed in Penalty Group 3
13	or 4;
14	Granulated opium;
15	Hydrocodone not listed in Penalty Group 3;
16	Hydromorphone;
17	Metopon;
18	Morphine not listed in Penalty Group 3;
19	Opium extracts;
20	Opium fluid extracts;
21	Oripavine;
22	Oxycodone;
23	Oxymorphone;
24	Powdered opium;
25	Raw opium;
26	Thebaine; and
27	Tincture of opium;

S.B. No. 1646 1 (B) a salt, compound, isomer, derivative, or preparation of a substance that is chemically equivalent or 2 3 identical to a substance described by Paragraph (A), other than the isoquinoline alkaloids of opium; 4 5 (C) Opium poppy and poppy straw; 6 (D) Cocaine, including: 7 (i) its salts, its optical, position, and 8 geometric isomers, and the salts of those isomers; 9 (ii) coca leaves and a salt, compound, 10 derivative, or preparation of coca leaves; and (iii) a salt, compound, derivative, 11 or preparation of a salt, compound, or derivative that is chemically 12 equivalent or identical to a substance described by Subparagraph 13 14 (i) or (ii), other than decocainized coca leaves or extractions of 15 coca leaves that do not contain cocaine or ecgonine; and 16 (E) concentrate of poppy straw, meaning the crude 17 extract of poppy straw in liquid, solid, or powder form that contains the phenanthrine alkaloids of the opium poppy; 18 the following opiates and opioids, including their 19 (4) isomers, esters, ethers, salts, and salts of isomers, if the 20 21 existence of these isomers, esters, ethers, and salts is possible within the specific chemical designation: 22 23 (A) [Acetyl-alpha-methylfentanyl 24 (N-[1-(1-methyl-2- phenethyl)-4-piperidinyl]-N-phenylacetamide); 25 [Alpha-methylthiofentanyl (N-[1-methyl-2-(2-26 thienyl)ethyl=4-piperidinyl]=N-phenylpropanamide); 27 Alphaprodine;

1	Anileridine;
2	[Beta=hydroxyfentanyl (N=[1=(2=hydroxy=2=
3	<pre>phenethyl)=4-piperidinyl] -N-phenylpropanamide);</pre>
4	<pre>[Beta=hydroxy=3=methylfentanyl;]</pre>
5	Bezitramide;
6	[Carfentanil;]
7	Dihydrocodeine not listed in Penalty Group 3 or 4;
8	Diphenoxylate not listed in Penalty Group 3 or 4;
9	Fentanyl [or alpha-methylfentanyl, or any other
10	<pre>derivative of Fentanyl];</pre>
11	Isomethadone;
12	Levomethorphan;
13	Levorphanol;
14	Metazocine;
15	Methadone;
16	Methadone-Intermediate, 4-cyano-2-dimethylamino-
17	4, 4-diphenyl butane;
18	[3-methylfentanyl(N-[3-methyl-1-(2-phenylethyl)-
19	4-piperidyl]-N-phenylpropanamide);
20	[3-methylthiofentanyl(N-[3-methyl-1-(2-thienyl)
21	<pre>ethyl=4-piperidinyl]-N-phenylpropanamide);</pre>
22	Moramide-Intermediate, 2-methyl-3-morpholino-1,
23	1-diphenyl-propane-carboxylic acid;
24	Para-fluorofentanyl(N-(4-fluorophenyl)-N-1-(2-
25	<pre>phenylethyl)-4-piperidinylpropanamide);</pre>
26	PEPAP (1-(2-phenethyl)-4-phenyl-4-
27	<pre>acetoxypiperidine);</pre>

	S.B. No. 1646
1	Pethidine (Meperidine);
2	Pethidine-Intermediate-A, 4-cyano-1-methyl-4-
3	phenylpiperidine;
4	Pethidine-Intermediate-B, ethyl-4-
5	phenylpiperidine-4 carboxylate;
6	Pethidine-Intermediate-C, 1-methyl-4-
7	phenylpiperidine-4-carboxylic acid;
8	Phenazocine;
9	Piminodine;
10	Racemethorphan; and
11	Racemorphan[;
12	[Remifentanil; and
13	[Thiofentanyl(N-phenyl-N-[1-(2-thienyl)ethyl-4-
14	<pre>piperidinyl]-propanamide)]; and</pre>
15	(B) a substance related to fentanyl, unless
16	specifically excepted or controlled in another Penalty Group,
17	including any substance that is structurally related to fentanyl by
18	one or more of the following modifications:
19	(i) substitution on or replacement of the
20	phenethyl group, such as:
21	Alpha methyl fentanyl;
22	Benzyl fentanyl; and
23	Thionyl fentanyl (HCl);
24	(ii) substitution on the piperidine ring,
25	such as:
26	<pre>Carfentanil;</pre>
27	3-Fluorofentanvl (HCl); and

4-Phenyl fentanyl (HCl); 1 2 (iii) substitution on the aniline ring, 3 such as: 4 para-Fluorofentanyl; 5 meta-Fluorofentanyl; and 6 ortho-Fluorofentanyl; and 7 (iv) substitution or replacement of the 8 N-propionyl group, such as: Acetyl fentanyl (HCL); 9 Benzodioxole fentanyl; and 10 Furanyl fentanyl (HCl); 11 Flunitrazepam (trade or other name: Rohypnol); 12 (5) Methamphetamine, including its salts, optical 13 (6) isomers, and salts of optical isomers; 14 15 (7) Phenylacetone and methylamine, if possessed 16 together with intent to manufacture methamphetamine; Phencyclidine, including its salts; 17 (8) Gamma hydroxybutyric acid (some trade or other 18 (9) names: gamma hydroxybutyrate, GHB), including its salts; 19 20 (10) Ketamine; 21 (11)Phenazepam; 22 (12) U-47700; 23 (13)AH-7921; 24 (14) ADB-FUBINACA; 25 (15)AMB-FUBINACA; and (16) MDMB-CHMICA. 26 SECTION 2. Section 481.103(a), Health and Safety Code, is 27

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amended to read as follows:
 1
 2
          (a)
              Penalty Group 2 consists of:
 3
               (1) any quantity of the following hallucinogenic
   substances, their salts, isomers, and salts of isomers, unless
 4
 5
   specifically excepted, if the existence of these salts, isomers,
    and salts of isomers is possible within the specific chemical
 6
 7
   designation:
8
                    5-(2-aminopropyl)benzofuran (5-APB);
 9
                    6-(2-aminopropyl)benzofuran (6-APB);
10
                    5-(2-aminopropyl)-2,3-dihydrobenzofuran
11
    (5-APDB);
12
                    6-(2-aminopropyl)-2,3-dihydrobenzofuran
    (6-APDB);
13
14
                    5-(2-aminopropyl) indole (5-IT, 5-API);
15
                    6-(2-aminopropyl)indole (6-IT,6-API);
16
                    1-(benzofuran-5-yl)-N-methylpropan-2-amine
17
    (5-MAPB);
                    1-(benzofuran-6-yl)-N-methylpropan-2-amine
18
19
    (6-MAPB);
20
                    Benzothiophenylcyclohexylpiperidine (BTCP);
21
                    8-bromo-alpha-methyl-benzo[1,2-b:4,5-b']difuran-
   4-ethanamine (trade or other name: Bromo-DragonFLY);
22
23
                    Desoxypipradrol (2-benzhydrylpiperidine);
24
                    2, 5-dimethoxyamphetamine (some trade or other
25
   names:
            2, 5-dimethoxy-alpha-methylphenethylamine; 2, 5-DMA);
26
                    Diphenylprolinol
                                          (diphenyl(pyrrolidin-2-yl)
27
   methanol, D2PM);
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S.B. No. 1646 1 Dronabinol (synthetic) in oil sesame and encapsulated in a soft gelatin capsule in a U.S. Food and Drug 2 3 Administration approved drug product (some trade or other names for Dronabinol: (a6aR-trans)-6a,7,8,10a-tetrahydro-4 6,6, 9-5 trimethyl-3-pentyl-6H- dibenzo [b,d]pyran-1-ol or (-)-delta-9-(trans) - tetrahydrocannabinol); 6 7 Ethylamine Analog of Phencyclidine (some trade or 8 other names: N-ethyl-1-phenylcyclohexylamine, (1 phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, 9 10 cyclohexamine, PCE); 2-ethylamino-2-(3-methoxyphenyl)cyclohexanone 11 12 (trade or other name: methoxetamine); Ibogaine (some trade or other names: 7-Ethyl-6, 13 6, beta 7, 8, 9, 10, 12, 13-octahydro-2-methoxy-6, 9-methano-5H-14 pyrido [1', 2':1, 2] azepino [5, 4-b] indole; tabernanthe iboga.); 15 5-iodo-2-aminoindane (5-IAI); 16 17 Mescaline; 5-methoxy-3, 4-methylenedioxy amphetamine; 18 19 4-methoxyamphetamine (some trade or other 20 names: 4-methoxy-alpha-methylphenethylamine; paramethoxyamphetamine; PMA); 21 4-methoxymethamphetamine (PMMA); 22 23 2-(2-methoxyphenyl)-2-(methylamino)cyclohexanone 24 (some trade and other names: 2-MeO-ketamine; methoxyketamine); 25 1-methyl- 4-phenyl-4-propionoxypiperidine (MPPP, 26 PPMP); 27 4-methyl-2, 5-dimethoxyamphetamine (some trade

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S.B. No. 1646
 1
   and
             other
                         names: 4-methyl-2,
                                                    5-dimethoxy-alpha-
   methylphenethylamine; "DOM"; "STP");
2
 3
                     3,4-methylenedioxy methamphetamine (MDMA, MDM);
 4
                     3,4-methylenedioxy amphetamine;
                                        N-ethylamphetamine
5
                     3,4-methylenedioxy
                                                                  (Also
   known as N-ethyl MDA);
6
7
                     5,6-methylenedioxy-2-aminoindane (MDAI);
8
                     Nabilone (Another name for nabilone: (+)-trans-
   3-(1,1-dimethylheptyl)- 6,6a, 7,8,10,10a-hexahydro-1-hydroxy- 6,
9
   6-dimethyl-9H-dibenzo[b,d] pyran-9-one;
10
11
                     N-benzylpiperazine
                                          (some
                                                   trade
                                                                 other
                                                            or
12
   names: BZP; 1-benzylpiperazine);
                     N-ethyl-3-piperidyl benzilate;
13
14
                     N-hydroxy-3,4-methylenedioxyamphetamine
                                                                  (Also
15
   known as N-hydroxy MDA);
16
                     4-methylaminorex;
17
                     N-methyl-3-piperidyl benzilate;
                     Parahexyl (some trade or other names: 3-Hexyl-1-
18
   hydroxy-7, 8, 9, 10-tetrahydro-6, 6, 9-trimethyl-6H-dibenzo [b, d]
19
20
   pyran; Synhexyl);
21
                     1-Phenylcyclohexylamine;
                     1-Piperidinocyclohexanecarbonitrile (PCC);
2.2
                     Pyrrolidine Analog of Phencyclidine (some trade
23
24
   or other names:
                    1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP);
25
                     Tetrahydrocannabinols, other than marihuana, and
26
   synthetic equivalents of the substances contained in the plant, or
    in the resinous extractives of Cannabis, or synthetic substances,
27
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S.B. No. 1646 1 derivatives, and their isomers with similar chemical structure, [and pharmacological activity] such as: 2 3 delta-9-tetrahydrocannabinolic acid (some trade name or other name: THCA-A and THCA-B); 4 5 delta-9-tetrahydrocannabinol, and its 6 optical isomers; 7 delta-1 cis or trans tetrahydrocannabinol, 8 and their optical isomers; 9 delta-6 cis or trans tetrahydrocannabinol, 10 and their optical isomers; delta-3, 11 4 cis trans or tetrahydrocannabinol, and its optical isomers; or 12 compounds of these structures, regardless of 13 14 numerical designation of atomic positions, since nomenclature of 15 these substances is not internationally standardized; 16 Thiophene Analog of Phencyclidine (some trade or 17 other names: 1-[1-(2-thienyl) cyclohexyl] piperidine; 2-Thienyl Analog of Phencyclidine; TPCP, TCP); 18 19 1-pyrrolidine (some trade or other name: TCPy); 20 1-(3-trifluoromethylphenyl)piperazine (trade or 21 other name: TFMPP); and 3,4,5-trimethoxy amphetamine; 22 23 (2) Phenylacetone (some trade or other Phenyl-2-propanone; P2P, Benzymethyl ketone, methyl benzyl 24 names: 25 ketone); 26 (3) unless specifically excepted or unless listed in Penalty Group, a material, compound, mixture, 27 another

12

or

S.B. No. 1646 1 preparation that contains any quantity of the following substances having a potential for abuse associated with a depressant or 2 3 stimulant effect on the central nervous system: 4 Aminorex (some trade or other 5 names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5phenyl-2-oxazolamine); 6 7 Amphetamine, its salts, optical isomers, and 8 salts of optical isomers; Cathinone (some trade or other names: 2-amino-1-9 10 phenyl-1-propanone, alpha-aminopropiophenone, 2aminopropiophenone); 11 12 Etaqualone and its salts; Etorphine Hydrochloride; 13 14 Fenethylline and its salts; 15 Lisdexamfetamine, including its salts, isomers, and salts of isomers; 16 17 Mecloqualone and its salts; Methaqualone and its salts; 18 Methcathinone (some trade or other names: 19 2methylamino-propiophenone; 20 alpha-(methylamino)propriophenone; 21 2-(methylamino)-1-phenylpropan-1-one; alpha-Nmethylaminopropriophenone; monomethylpropion; ephedrone, 22 Nmethylcathinone; methylcathinone; AL-464; AL-422; AL-463; and UR 23 24 1431); 25 N-Ethylamphetamine, its salts, optical isomers, 26 and salts of optical isomers; and 27 N,N-dimethylamphetamine (some trade or other

S.B. No. 1646 1 names: N,N,alpha-trimethylbenzeneethanamine; N,N,alpha-trimethylphenethylamine), its salts, optical isomers, 2 3 and salts of optical isomers; 4 (4) any compound structurally derived from 5 2-aminopropanal by substitution at the 1-position with any monocyclic or fused-polycyclic ring system, including: 6 7 (A) compounds further modified by: 8 (i) substitution in the ring system to any extent (including alkyl, alkoxy, alkylenedioxy, haloalkyl, or 9 halide substituents), whether or not further substituted in the 10 ring system by other substituents; 11 12 (ii) substitution at the 3-position with an alkyl substituent; or 13 14 (iii) substitution at the 2-amino nitrogen 15 atom with alkyl, benzyl, dialkyl, or methoxybenzyl groups, or inclusion of the 2-amino nitrogen atom in a cyclic structure; and 16 17 (B) by example, compounds such as: 4-Methylmethcathinone 18 (Also known as 19 Mephedrone); 3,4-Dimethylmethcathinone 20 (Also known as 3,4-DMMC); 21 3-Fluoromethcathinone (Also known as 3-FMC); 2.2 23 4-Fluoromethcathinone (Also known as 24 Flephedrone); 25 3,4-Methylenedioxy-N-methylcathinone (Also 26 known as Methylone); 27 3,4-Methylenedioxypyrovalerone (Also known

1 as MDPV); 2 alpha-Pyrrolidinopentiophenone (Also known 3 as alpha-PVP); 4 Naphthylpyrovalerone (Also known as 5 Naphyrone); 6 alpha-Methylamino-valerophenone (Also known 7 as Pentedrone); 8 beta-Keto-N-methylbenzodioxolylpropylamine (Also known as Butylone); 9 10 beta-Keto-N-methylbenzodioxolylpentanamine (Also known as Pentylone); 11 beta-Keto-Ethylbenzodioxolylbutanamine 12 (Also known as Eutylone); and 13 14 3,4-methylenedioxy-N-ethylcathinone (Also 15 known as Ethylone); 16 (5) any compound structurally derived from tryptamine 17 (3-(2-aminoethyl)indole) or a ring-hydroxy tryptamine: (A) by modification in any of the following ways: 18 (i) by substitution at the amine nitrogen 19 atom of the sidechain to any extent with alkyl or alkenyl groups or 20 by inclusion of the amine nitrogen atom of the side chain (and no 21 22 other atoms of the side chain) in a cyclic structure; 23 (ii) by substitution at the carbon atom 24 adjacent to the nitrogen atom of the side chain (alpha-position) 25 with an alkyl or alkenyl group; (iii) by substitution in the 6-membered 26 ring to any extent with alkyl, alkoxy, haloalkyl, thioaklyl, 27

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S.B. No. 1646
   alkylenedioxy, or halide substituents; or
 1
 2
                          (iv) by substitution at the 2-position of
 3
   the tryptamine ring system with an alkyl substituent; and
 4
                     (B)
                         including:
 5
                          (i) ethers and esters of the controlled
   substances listed in this subdivision; and
 6
 7
                          (ii) by example, compounds such as:
8
                               alpha-ethyltryptamine;
 9
                               alpha-methyltryptamine;
                               Bufotenine (some trade and other names:
10
   3-(beta-Dimethylaminoethyl)-5-hydroxyindole;
11
   3-(2-dimethylaminoethyl)- 5- indolol; N, N-dimethylserotonin;
12
   5-hydroxy-N, N- dimethyltryptamine; mappine);
13
14
                               Diethyltryptamine
                                                   (some
                                                           trade
                                                                   and
15
   other names: N, N-Diethyltryptamine, DET);
16
                               Dimethyltryptamine
                                                    (trade
                                                                 other
                                                            or
17
   name: DMT);
                               5-methoxy-N,
                                              N-diisopropyltryptamine
18
    (5-MeO-DiPT);
19
20
                               O-Acetylpsilocin (Trade or other name:
   4-Aco-DMT);
21
                               Psilocin; and
2.2
23
                               Psilocybin;
24
               (6)
                    2,5-Dimethoxyphenethylamine
                                                 and
                                                        any
                                                             compound
   structurally derived
                            from 2,5-Dimethoxyphenethylamine
25
                                                                    by
   substitution at the 4-position of the phenyl ring to any extent
26
    (including alkyl, alkoxy, alkylenedioxy, haloalkyl, or halide
27
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16
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S.B. No. 1646
 1
    substituents), including, by example, compounds such as:
                     4-Bromo-2,5-dimethoxyphenethylamine
 2
                                                            (trade
                                                                     or
 3
    other name: 2C-B);
 4
                     4-Chloro-2,5-dimethoxyphenethylamine
                                                             (trade
                                                                     or
 5
    other name: 2C-C);
                     2,5-Dimethoxy-4-methylphenethylamine
 6
                                                             (trade
                                                                     or
 7
    other name: 2C-D);
8
                     4-Ethyl-2,5-dimethoxyphenethylamine
                                                            (trade
                                                                     or
    other name: 2C-E);
 9
10
                     4-Iodo-2,5-dimethoxyphenethylamine
                                                            (trade
                                                                     or
   other name: 2C-I);
11
12
                     2,5-Dimethoxy-4-nitrophenethylamine
                                                            (trade
                                                                     or
    other name: 2C-N);
13
14
                     2,5-Dimethoxy-4-(n)-propylphenethylamine (trade
15
    or other name: 2C-P);
                     4-Ethylthio-2,5-dimethoxyphenethylamine
16
                                                                 (trade
17
    or other name: 2C-T-2);
                     4-Isopropylthio-2,5-dimethoxyphenethylamine
18
19
    (trade or other name: 2C-T-4); and
                     2,5-Dimethoxy-4-(n)-propylthiophenethylamine
20
21
    (trade or other name: 2C-T-7); and
                     2,5-Dimethoxyamphetamine
22
                (7)
                                                 and
                                                               compound
                                                        any
    structurally derived from 2,5-Dimethoxyamphetamine by substitution
23
24
    at the 4-position of the phenyl ring to any extent (including alkyl,
    alkoxy, alkylenedioxy, haloalkyl, or
25
                                                halide substituents),
26
    including, by example, compounds such as:
27
                     4-Ethylthio-2,5-dimethoxyamphetamine (trade
                                                                     or
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17
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1
    other name: Aleph-2);
 2
                     4-Isopropylthio-2,5-dimethoxyamphetamine (trade
 3
    or other name: Aleph-4);
 4
                     4-Bromo-2,5-dimethoxyamphetamine (trade or other
 5
    name: DOB);
                     4-Chloro-2,5-dimethoxyamphetamine
 6
                                                           (trade
                                                                    or
 7
    other name: DOC);
                     2,5-Dimethoxy-4-ethylamphetamine (trade or other
8
    name: DOET);
 9
10
                     4-Iodo-2,5-dimethoxyamphetamine (trade or other
11
   name: DOI);
12
                     2,5-Dimethoxy-4-methylamphetamine
                                                          (trade
                                                                    or
13
    other name: DOM);
14
                     2,5-Dimethoxy-4-nitroamphetamine (trade or other
15
   name: DON);
16
                     4-Isopropyl-2,5-dimethoxyamphetamine (trade
                                                                    or
17
    other name: DOIP); and
                     2,5-Dimethoxy-4-(n)-propylamphetamine (trade
18
                                                                    or
19
    other name: DOPR).
          SECTION 3. Section 481.1031, Health and Safety Code, is
20
    amended to read as follows:
21
          Sec. 481.1031. PENALTY GROUP 2-A. (a)
                                                 In this section:
22
               (1) "Core component" is
23
                                                  of
                                             one
                                                      the
                                                            following:
24
    azaindole,
                 beta carboline, benzimidazole,
                                                       benzothiazole,
    carbazole, gamma carboline, imidazole, indane, indazole, indene,
25
26
    indole, <u>naphthalene</u>, pyrazole, pyrazolopyridine, pyridine, or
   pyrrole.
27
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1 (2) "Group A component" is one of the following: benzylpiperazine, benzyl, cumene 2 adamantane, [benzene], 3 cycloalkylmethyl, cyclopropylmethyl, isoquinoline, methylpiperazine, naphthalene, pyrrolidine, phenyl, quinoline, 4 5 tetrahydronaphthalene, tetramethylcyclopropane, amino oxobutane, amino methyl ovobutane, amino dimethyl oxobutane, amino phenyl 6 oxopropane, amino oxopentane, ethoxy oxobutane, ethoxy methyl 7 oxobutane, ethoxy dimethyl oxobutane, ethoxy phenyl oxopropane, 8 ethoxy oxopentane, methoxy oxobutane, methoxy methyl oxobutane 9 10 [methyl methoxy oxobutane], methoxy dimethyl oxobutane, methoxy phenyl oxopropane, methoxy oxopentane, or any substituted [an] 11 12 amino acid.

S.B. No. 1646

(3) "Link component" 13 is one of the following 14 functional groups: carboxamide, carboxylate, hydrazide, 15 methanone (ketone), ethanone, methanediyl (methylene bridge), or 16 methine.

(b) Penalty Group 2-A consists of any material, compound, mixture, or preparation that contains any quantity of a natural or synthetic chemical substance, including its salts, isomers, and calts of isomers, listed by name in this subsection or contained within one of the structural classes defined in this subsection:

22

(1) WIN-55,212-2;

(2) Cyclohexylphenol: any compound structurally
derived from 2-(3-hydroxycyclohexyl)phenol by substitution at the
5-position of the phenolic ring, (N-methylpiperidin-2-yl)alkyl,
(4-tetrahydropyran)alkyl, or 2-(4-morpholinyl)alkyl, whether or
not substituted in the cyclohexyl ring to any extent, including:

1 JWH-337; 2 JWH-344; CP-55,940; 3 CP-47,497; and 4 5 analogues of CP-47,497; 6 (3) Cannabinol derivatives, except where contained in 7 marihuana, including tetrahydro derivatives of cannabinol and 8 3-alkyl homologues of cannabinol or of its tetrahydro derivatives, 9 such as: Nabilone; 10 HU-210; and 11 12 HU-211; (4) Tetramethylcyclopropyl thiazole: any compound 13 14 structurally derived from 2,2,3,3-tetramethyl-N-(thiazol-15 2-ylidene)cyclopropanecarboxamide by substitution at the nitrogen atom of the thiazole ring, whether or not further substituted in the 16 17 thiazole ring to any extent, whether or not substituted in the tetramethylcyclopropyl ring to any extent, such as [including]: 18 A-836,339; 19 20 (5) any compound containing a core component 21 substituted at the 1-position to any extent, and substituted at the 3-position with a link component attached to a group A component, 22 whether or not the core component or group A component are further 23 24 substituted to any extent, <u>such as</u> [including]: 25 Naphthoylindane; Naphthoylindazole (THJ-018); 26 Naphthyl methyl indene (JWH-171); 27

Naphthoylindole (JWH-018); 1 2 Quinolinoyl pyrazole carboxylate (Quinolinyl 3 fluoropentyl fluorophenyl pyrazole carboxylate); 4 Naphthoyl pyrazolopyridine; and 5 Naphthoylpyrrole (JWH-030); 6 (6) any compound containing a core component substituted at the 1-position to any extent, and substituted at the 7 8 2-position with a link component attached to a group A component, whether or not the core component or group A component are further 9 substituted to any extent, <u>such as</u> [including]: 10 Naphthoylbenzimidazole (JWH-018 Benzimidazole); 11 12 and Naphthoylimidazole; 13 14 (7)any compound containing a core component 15 substituted at the 3-position to any extent, and substituted at the 2-position with a link component attached to a group A component, 16 17 whether or not the core component or group A component are further substituted to any extent, such as [including]: 18 Naphthoyl benzothiazole; [and] 19 20 (8) containing a any compound core component 21 substituted at the 9-position to any extent, and substituted at the 3-position with a link component attached to a group A component, 22 23 whether or not the core component or group A component are further 24 substituted to any extent, including: 25 Naphthoylcarbazole (EG-018); and 26 Synthetic chemical compounds with a carbazole core structure, regardless of numerical designation of atomic 27

1 positions, since this core structure is symmetrical; 2 (9) any compound containing a core component 3 substituted at the 1-position to any extent and substituted at the 5-position with a link component attached to a group A component, 4 5 regardless of whether the core component or group A component are 6 further substituted to any extent, such as: 7 5-fluoro-3,5-ADB-PFUPPYCA (5f-AB-PFUPPYCA); (10) any compound containing a core component 8 substituted at the 2-position to any extent and substituted at the 9 10 3-position with a link component attached to a group A component, regardless of whether the core component or group A component are 11 12 further substituted to any extent, such as: AB-CHMINACA (2H) indazole; and 13 (11) any compound containing a core component 14 15 substituted at the 5-position to any extent and substituted at the 2-position with a group A component, regardless of whether the core 16 17 component or group A component are further substituted to any extent, such as: 18 19 Cumyl-PeGACLONE. (c) In this section, a synthetic chemical substance 20 21 analogue is: 22 (1) any substance, unless specifically excepted, that has two of the three components as defined by Subsection (a) and a 23 24 numerical core-link position that is listed in Subsection (b)(5), (6), (7), (8), (9), or (10); 25 26 (2) any substance, unless specifically excepted, that 27 has a core, link, and group A component as defined by Subsection (a)

1 but does not have a numerical core-link position that is listed in 2 Subsection (b)(5), (6), (7), (8), (9), or (10); or

3 (3) any substance, unless specifically excepted, that 4 has a core component or group A component as defined by Subsection 5 (a) and a numerical core-group A position that is listed in 6 Subsection (b)(11).

SECTION 4. The change in law made by this Act applies only 7 8 to an offense committed on or after the effective date of this Act. An offense committed before the effective date of this Act is 9 governed by the law in effect on the date the offense was committed, 10 and the former law is continued in effect for that purpose. For 11 purposes of this section, an offense was committed before the 12 effective date of this Act if any element of the offense occurred 13 14 before that date.

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SECTION 5. This Act takes effect September 1, 2019.