```
1/13/20 7:29
```

```
20102765D
 1
                                           SENATE BILL NO. 538
 2
                                            Offered January 8, 2020
 3
                                           Prefiled January 7, 2020
    A BILL to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Drug Control Act;
 5
        controlled substances: Schedule I.
 6
                                              Patron—Newman
 7
 8
                                 Referred to Committee on Education and Health
 9
10
        Be it enacted by the General Assembly of Virginia:
    1. That § 54.1-3446 of the Code of Virginia is amended and reenacted as follows:
11
        § 54.1-3446. Schedule I.
12
13
        The controlled substances listed in this section are included in Schedule I:
14
        1. Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers,
15
    esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers
    and salts is possible within the specific chemical designation:
16
        1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP);
17
        1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP);
18
19
        2-methoxy-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide
                                                                         (other
                                                                                            Methoxyacetyl
                                                                                   name:
20
    fentanvl):
21
        3.4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzamide (other name: U-47700);
22
        3,4-dichloro-N-{[1-(dimethylamino)cyclohexyl]methyl]benzamide (other name: AH-7921);
23
        Acetyl fentanyl (other name: desmethyl fentanyl);
24
        Acetylmethadol;
25
        Allylprodine;
        Alphacetylmethadol (except levo-alphacetylmethadol, also known as levo-alpha-acetylmethadol,
26
27
    levomethadyl acetate, or LAAM);
28
        Alphameprodine:
29
        Alphamethadol;
30
        Benzethidine;
31
        Betacetylmethadol;
        Betameprodine;
32
33
        Betamethadol;
34
        Betaprodine;
35
        Clonitazene:
36
        Dextromoramide:
37
        Diampromide;
        Diethylthiambutene;
38
39
        Difenoxin;
        Dimenoxadol:
40
41
        Dimepheptanol:
        Dimethylthiambutene;
42
43
        Dioxaphetylbutyrate;
        Dipipanone:
44
        Ethylmethylthiambutene;
45
46
        Etonitazene:
47
        Etoxeridine:
48
        Furethidine:
49
        Hydroxypethidine;
        Ketobemidone:
50
51
        Levomoramide;
52
        Levophenacylmorphan;
53
        Morpheridine:
54
        MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
55
        N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (other name: Cyclopropyl fentanyl);
        N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (other name: Tetrahydrofuranyl
56
57
        N-[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]-N-phenylpropanamide (other name:
58
```

SB538 2 of 8

120

```
59
         alpha-methylthiofentanyl);
 60
             N-[1-(1-methyl-2-phenylethyl)-4-piperidyl]-N-phenylacetamide (other name:
 61
         acetyl-alpha-methylfentanyl);
 62
             N-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl}-N-phenylpropanamide (other name:
 63
         beta-hydroxythiofentanyl);
             N-[1-(2-hydroxy-2-phenyl)ethyl-4-piperidyl]-N-phenylpropanamide (other name:
 64
 65
         beta-hydroxyfentanyl);
 66
             N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide (other names:
         1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine, alpha-methylfentanyl);
 67
             N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other names: 2-fluorofentanyl,
 68
 69
         ortho-fluorofentanyl):
 70
             N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: 3-fluorofentanyl);
             N-[3-methyl-1-(2-hydroxy-2-phenylethyl)4-piperidyl]-N-phenylpropanamide (other name:
 71
 72
         beta-hydroxy-3-methylfentanyl);
             N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (other name: 3-methylfentanyl);
 73
 74
             N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (other name:
 75
         3-methylthiofentanyl);
 76
             N-(4-fluorophenyl)-2-methyl-N-[1-(2-phenylethyl)-4-piperidinyl] -propanamide (other name:
 77
         para-fluoroisobutyryl fentanyl);
 78
             N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name:
 79
         para-fluorobutyrylfentanyl);
 80
             N-(4-fluorophenyl)-N-1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: para-fluorofentanyl);
 81
              Noracymethadol;
 82
             Norlevorphanol;
 83
             Normethadone;
 84
             Norpipanone:
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-furancarboxamide (other name: Furanyl fentanyl);
 85
 86
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-propenamide (other name: Acryl fentanyl);
 87
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: butyryl fentanyl);
 88
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide (other name: Pentanoyl fentanyl);
 89
             N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide (other name: thiofentanyl);
 90
             Phenadoxone:
 91
             Phenampromide;
 92
             Phenomorphan;
 93
             Phenoperidine:
 94
             Piritramide:
 95
             Proheptazine;
 96
             Properidine;
 97
             Propiram;
 98
             Racemoramide:
 99
             Tilidine:
100
             Trimeperidine:
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-1,3-benzodioxole-5 -carboxamide (other name:
101
102
         Benzodioxole fentanyl):
              3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methylbenzamide (other name: U-49900);
103
104
             2-(2,4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methylacetamide (other name: U-48800);
105
             2-(3.4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methylacetamide (other name: U-51754);
             N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide (other name: Ocfentanil);  
    <math>N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: N-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxyphenyl)-N-[1-(4-methoxypheny
106
107
108
         4-methoxybutyrylfentanyl);
109
             N-phenyl-2-methyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: Isobutyryl fentanyl);
110
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-cyclopentanecarboxamide (other name: Cyclopentyl
111
         fentanyl);
             N-phenyl-N-(1-methyl-4-piperidinyl)-propanamide (other name: N-methyl norfentanyl);
112
             N-[2-(dimethylamino)cyclohexyl]-N-methyl-1,3-benzodioxole-5-carboxamide (other names:
113
114
         3.4-methylenedioxy U-47700 or 3.4-MDO-U-47700);
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-butenamide (other name: Crotonyl fentanyl);
115
116
             N-phenyl-N-[4-phenyl-1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: 4-phenylfentanyl);
             N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-benzamide (other names: Phenyl fentanyl, Benzoyl
117
118
             N-[2-(dimethylamino)cyclohexyl]-N-phenylfuran-2-carboxamide (other name: Furanyl UF-17);
119
```

*N-[2-(dimethylamino)cyclohexyl]-N-phenylpropionamide (other name: UF-17);* 

```
121
        3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-isopropyl-benzamide (other name: Isopropyl
122
     U-47700).
```

123 2. Any of the following opium derivatives, their salts, isomers and salts of isomers, unless 124 specifically excepted, whenever the existence of these salts, isomers and salts of isomers is possible 125 within the specific chemical designation:

126 Acetorphine:

127 Acetyldihydrocodeine;

128 Benzylmorphine;

- 129 Codeine methylbromide;
- 130 Codeine-N-Oxide;
- 131 Cyprenorphine;
- 132 Desomorphine;
- 133 Dihydromorphine;
- 134 Drotebanol;
- 135 Etorphine:
- 136 Heroin;
- 137 Hydromorphinol;
- 138 Methyldesorphine:
- 139 Methyldihydromorphine;
- 140 Morphine methylbromide;
- 141 Morphine methylsulfonate:
- 142 Morphine-N-Oxide;
- 143 Myrophine;
- 144 Nicocodeine:
- 145 Nicomorphine;
- 146 Normorphine;
- 147 Pholcodine:
- 148 Thebacon.

149

150

151

152

153

154

3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or preparation, which contains any quantity of the following hallucinogenic substances, or which contains any of 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 (for purposes of this subdivision only, the term "isomer" includes the optical, position, and geometric isomers):

Alpha-ethyltryptamine (some trade or other names: Monase; a-ethyl-1H-indole-3-ethanamine; 3-2-aminobutyl] indole; a-ET; AET);

155

156 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names: 157 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);

158 3.4-methylenedioxy amphetamine:

- 159 5-methoxy-3,4-methylenedioxy amphetamine;
- 160 3,4,5-trimethoxy amphetamine;
- Alpha-methyltryptamine (other name: AMT); 161
- 162 Bufotenine;
- Diethyltryptamine; 163
- Dimethyltryptamine; 164
- 4-methyl-2,5-dimethoxyamphetamine; 165
- 166 2,5-dimethoxy-4-ethylamphetamine (DOET);
- 167 4-fluoro-N-ethylamphetamine;
- 168 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
- 169 Ibogaine;
- **170** 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
- 171 Lysergic acid diethylamide;
- 172 Mescaline;
- 173 Parahexyl (some trade other names: or
- 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenz o [b,d] pyran; Synhexyl); 174
- 175 Peyote;
- 176 N-ethyl-3-piperidyl benzilate;
- 177 N-methyl-3-piperidyl benzilate;
- 178 Psilocybin;
- 179 Psilocyn;
- 180 Salvinorin A:
- 181 Tetrahydrocannabinols, except as present in (i) industrial hemp, as defined in § 3.2-4112, that is

SB538 4 of 8

182 possessed by a person registered pursuant to subsection A of § 3.2-4115 or his agent; (ii) a hemp product, as defined in § 3.2-4112, containing a tetrahydrocannabinol concentration of no greater than 0.3 183 184 percent that is derived from industrial hemp, as defined in § 3.2-4112, that is grown, dealt, or processed 185 in compliance with state or federal law; (iii) marijuana; or (iv) dronabinol in sesame oil and 186 encapsulated in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug 187 Administration: 188 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish); 189 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine; 190 191 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts 192 and salts of isomers; 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl-alpha-methyl-3,4 193 194 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA); 195 N-hydroxy-3,4-methylenedioxyamphetamine (some other names: N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA); 196 197 4-bromo-2,5-dimethoxyamphetamine (some trade other names: 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA); 198 199 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine; 200 paramethoxyamphetamine; PMA); Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine, 201 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE); 202 203 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy, PHP); 204 205 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl) -cyclohexyl]-piperidine, 2-thienyl analog of phencyclidine, TPCP, TCP); 206 1-1-(2-thienyl)cyclohexyl]pyrrolidine (other name: TCPy); 207 208 3,4-methylenedioxypyrovalerone (other name: MDPV); 209 4-methylmethcathinone (other names: mephedrone, 4-MMC); 210 3.4-methylenedioxymethcathinone (other name: methylone); Naphthylpyrovalerone (other name: naphyrone); 211 212 4-fluoromethcathinone (other name: flephedrone, 4-FMC); 213 4-methoxymethcathinone (other names: methodrone; bk-PMMA); 214 Ethcathinone (other name: N-ethylcathinone); 215 3,4-methylenedioxyethcathinone (other name: ethylone); 216 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone); 217 N,N-dimethylcathinone (other name: metamfepramone); 218 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP); 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP); 219 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP); 220 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP); 221 222 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI); 223 3-fluoromethcathinone (other name: 3-FMC); 224 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E); 225 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I); 226 4-Methylethcathinone (other name: 4-MEC); 227 4-Ethylmethcathinone (other name: 4-EMC); 228 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT); 229 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP); 230 Alpha-methylamino-butyrophenone (other name: Buphedrone); 231 Alpha-methylamino-valerophenone (other name: Pentedrone); 3,4-Dimethylmethcathinone (other name: 3.4-DMMC); 232 233 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP); 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I, 234 25I-NBOMe, 2C-I-NBOMe); 235 236 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE); 237 4-Fluoromethamphetamine (other name: 4-FMA); 238 4-Fluoroamphetamine (other name: 4-FA); 239 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D); 240 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C); 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2); 241

2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4); 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H); 243

242

```
244
         2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
245
         2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
246
         (2-aminopropyl)benzofuran (other name: APB);
247
         (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
         4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
248
249
     2C-C-NBOMe, 25C-NBOMe, 25C);
250
         4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:
251
     2C-B-NBOMe, 25B-NBOMe, 25B);
252
         Acetoxydimethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
253
         Benocyclidine (other names: BCP, BTCP);
254
         Alpha-pyrrolidinobutiophenone (other name: alpha-PBP);
255
         3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
256
         4-bromomethcathinone (other name: 4-BMC);
257
         4-chloromethcathinone (other name: 4-CMC);
         4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
258
259
         Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
260
         Alpha-Pyrrolidinoheptiophenone (other name: PV8);
261
         5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
262
         Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
263
         Beta-keto-4-bromo-2,5-dimethoxyphenethylamine (other name: bk-2C-B);
264
         1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
265
         1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
266
         1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
267
         4-Chloroethcathinone (other name: 4-CEC);
268
         3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
269
         1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
270
         (2-Methylaminopropyl)benzofuran (other name: MAPB);
271
         1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)-1-pentanone
                                                                (other
                                                                                  N,N-Dimethylpentylone,
                                                                        names:
272
     Dipentylone);
273
         1-(4-methoxyphenyl)-2-(pyrrolidin-1-yl)octan-1-one (other name: 4-methoxy-PV9);
274
         3,4-tetramethylene-alpha-pyrrolidinovalerophenone (other name: TH-PVP);
275
         4-allyloxy-3,5-dimethoxyphenethylamine (other name: Allylescaline);
276
         4-Bromo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25B-NBOH);
277
         4-chloro-alpha-methylamino-valerophenone (other name: 4-chloropentedrone);
278
         4-chloro-alpha-Pyrrolidinovalerophenone (other name: 4-chloro-alpha-PVP);
279
         4-fluoro-alpha-Pyrrolidinoheptiophenone (other name: 4-fluoro-PV8);
280
         4-hydroxy-N,N-diisopropyltryptamine (other name: 4-OH-DIPT);
281
         4-methyl-alpha-ethylaminopentiophenone;
282
         4-methyl-alpha-Pyrrolidinohexiophenone (other name: MPHP);
         5-methoxy-N,N-dimethyltryptamine (other name: 5-MeO-DMT);
283
284
         5-methoxy-N-ethyl-N-isopropyltryptamine (other name: 5-MeO-EIPT);
285
         6-ethyl-6-nor-lysergic acid diethylamide (other name: ETH-LAD);
286
         6-allyl-6-nor-lysergic acid diethylamide (other name: AL-LAD);
287
         (N-methyl aminopropyl)-2,3-dihydrobenzofuran (other name: MAPDB);
288
         2-(methylamino)-2-phenyl-cyclohexanone (other name: Deschloroketamine);
289
         2-(ethylamino)-2-phenyl-cyclohexanone (other name: deschloro-N-ethyl-ketamine);
290
         2-methyl-1-(4-(methylthio)phenyl)-2-morpholinopropiophenone (other name: MMMP);
291
         Alpha-ethylaminohexanophenone (other name: N-ethylhexedrone);
292
         N-ethyl-1-(3-methoxyphenyl)cyclohexylamine (other name: 3-methoxy-PCE);
293
         4-fluoro-alpha-pyrrolidinohexiophenone (other name: 4-fluoro-alpha-PHP);
294
         N-ethyl-1,2-diphenylethylamine (other name: Ephenidine);
295
         2,5-dimethoxy-4-chloroamphetamine (other name: DOC);
296
         3,4-methylenedioxy-N-tert-butylcathinone;
297
        Alpha-pyrrolidinoisohexiophenone (other name: alpha-PiHP);
298
         1-[1-(3-hydroxyphenyl)cyclohexyl]piperidine (other name: 3-hydroxy PCP);
299
         4-acetyloxy-N,N-diallyltryptamine (other name: 4-AcO-DALT);
300
         4-hydroxy-N,N-methylisopropyltryptamine (other name: 4-hydroxy-MiPT);
301
         3,4-Methylenedioxy-alpha-pyrrolidinohexanophenone (other name: MDPHP);
302
         5-methoxy-N,N-dibutyltryptamine (other name: 5-methoxy-DBT);
         1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (other name: Eutylone, bk-EBDB);
303
```

1-(1,3-benzodioxol-5-yl)-2-(butylamino)-1-pentanone (other name: N-butylpentylone);

304

SB538 6 of 8

- 305 N-benzyl-3,4-dimethoxyamphetamine (other name: N-benzyl-3,4-DMA).
- 306 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture 307 or preparation which contains any quantity of the following substances having a depressant effect on the 308 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such 309 salts, isomers and salts of isomers is possible within the specific chemical designation:

310 Clonazolam:

311 Etizolam;

317

318 319

320

321

322

323

324

325

326

327

330

331

332

333 334

335

339

340

341

342

343 344

345

346 347 348

349

350

351

352

353

354

355

356 357

358 359

360

361

362 363

364 365

366

312 Flualprazolam;

Flubromazepam; 313 314

Flubromazolam;

315 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate); 316

Mecloqualone;

Methaqualone.

5. 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:

2-(3-fluorophenyl)-3-methylmorpholine (other name: 3-fluorophenmetrazine);

Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-phenyl-2-oxazolamine);

Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone, 2-aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;

Cis-4-methylaminorex (other name: cis-4.5-dihydro-4-methyl-5-phenyl-2-oxazolamine);

328 Ethylamphetamine: 329

Ethyl phenyl(piperidin-2-yl)acetate (other name: Ethylphenidate);

Fenethylline:

Methcathinone names: 2-(methylamino)-propiophenone; (some other alpha-(methylamino)-propiophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone; methylcathinone; AL-464; AL-422; AL-463 and UR 1432);

N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);

336 N,N-dimethylamphetamine (other names: N, N-alpha-trimethyl-benzeneethanamine, N, 337 N-alpha-trimethylphenethylamine); 338

Methyl 2-(4-fluorophenyl)-2-(2-piperidinyl)acetate (other name: 4-fluoromethylphenidate);

Isopropyl-2-phenyl-2-(2-piperidinyl)acetate (other name: Isopropylphenidate);

4-chloro-N,N-dimethylcathinone;

- *3,4-methylenedioxy-N-benzylcathinone (other name: BMDP).*
- 6. Any substance that contains one or more cannabimimetic agents or that contains their salts, isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation, and any preparation, mixture, or substance containing, or mixed or infused with, any detectable amount of one or more cannabimimetic agents.
- a. "Cannabimimetic agents" includes any substance that is within any of the following structural
- 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or alkenyl, whether or not substituted on the cyclohexyl ring to any extent;
- 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl)methane with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the naphthoyl or naphthyl ring to any extent;
- 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to
- 1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to
- 3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted in the indole ring to any extent, whether or not substituted on the phenyl ring to any extent;
- 3-cyclopropoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any
- 3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any

```
367
368
         N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring,
369
     whether or not further substituted on the indole ring to any extent, whether or not substituted on the
370
     adamantyl ring to any extent; and
371
         N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring,
372
     whether or not further substituted on the indazole ring to any extent, whether or not substituted on the
373
     adamantyl ring to any extent.
374
         b. The term "cannabimimetic agents" includes:
375
         5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);
376
         5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);
377
         5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);
378
         5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);
379
         1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);
380
         1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);
381
         1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);
382
         1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);
383
         1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);
384
         (6aR, 10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-ter
385
     ahydrobenzo[c]chromen-1-ol (other name: HU-210);
386
         1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);
387
         1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);
388
         1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);
389
         1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);
390
         1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);
391
         1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);
392
         1-((N-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);
393
         1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);
394
         1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);
395
                       (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone
         Pravadoline
                                                                                                   (other
396
     name: WIN 48,098);
397
         1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);
398
         1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);
399
         1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);
400
         1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other names: XLR-11,
401
     5-fluoro-UR-144);
402
         N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);
403
         N-adamantyl-1-pentylindazole-3-carboxamide (other names: AKB48, APINACA);
404
         1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);
405
         (8-quinolinyl)(1-pentylindol-3-yl)carboxylate (other name: PB-22);
406
         (8-quinolinyl)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);
407
         (8-quinolinyl)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);
408
         N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);
409
         N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name:
410
     AB-FUBINACA);
411
         1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);
412
         N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name:
     ADB-PINACA);
413
414
         N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide
                                                                                          (other
                                                                                                   name:
415
     AB-CHMINACA);
416
        N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:
417
     5-fluoro-AB-PINACA);
418
         N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxa mide (other
     names: ADB-CHMINACA, MAB-CHMINACA);
419
420
         Methyl-2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name:
421
     5-fluoro-AMB);
422
         1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);
423
         1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);
424
         1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name MAM-2201);
425
         N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-[(4-fluorophenyl)methyl]-1H-indazole-
                                                                                           3-carboxamide
```

Methyl 2-[1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name:

426

427

(other name: ADB-FUBINACA);

SB538 8 of 8

- **428** MDMB-FUBINACA);
- Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names: 5-fluoro-ADB, 5-Fluoro-MDMB-PINACA;
- 431 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoat e (other 432 names: AMB-FUBINACA, FUB-AMB);
- N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other name: FUB-AKB48);
- N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (other name: 5F-AKB48);
- N-(adamantanyl)-1-(5-chloropentyl) indazole-3-carboxamide (other name: 5-chloro-AKB48);
- Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate (other name: SDB-005);
- N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide (other name: 438 AB-CHMICA);
- 439 1-pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide (other name: SDB-006);
- Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate (other name: FUB-PB-22);
- Methyl N-[1-(cyclohexylmethyl)-1H-indole-3-carbonyl]valinate (other name: MMB-CHMICA);
- N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamid e (other name: 5-fluoro-ADB-PINACA);
- 1-(4-cyanobutyl)-N-(1-methyl-1-phenylethyl)-1H-indazole-3-carboxamide (other name: 4-cyano CUMYL-BUTINACA);
- 446 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name: 447 5-Fluoro-MDMB-PICA);
- 448 Ethyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other name: 449 EMB-FUBINACA);
- 450 Methyl 2-[1-4-fluorobutyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name: 451 4-fluoro-MDMB-BUTINACA).
- 2. That the provisions of this act may result in a net increase in periods of imprisonment or commitment. Pursuant to § 30-19.1:4 of the Code of Virginia, the estimated amount of the necessary appropriation is \$0 for periods of imprisonment in state adult correctional facilities and cannot be determined for periods of commitment to the custody of the Department of Juvenile
- 456 Justice.