

HOUSE BILL NO. 558

AMENDMENT IN THE NATURE OF A SUBSTITUTE

(Proposed by the House Committee on Health and Human Services

on _____)

(Patron Prior to Substitute—Delegate Wachsmann, Jr.)

A BILL to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Drug Control Act; Schedule I; penalties.

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:****§ 54.1-3446. Schedule I.**

The controlled substances listed in this section are included in Schedule I:

1. Any of the following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and ethers, unless specifically excepted, whenever the existence of these isomers, esters, ethers and salts is possible within the specific chemical designation:

1-{1-[1-(4-bromophenyl)ethyl]-4-piperidinyl}-1,3-dihydro-2H-benzimidazol-2-one (other name: Brorphine);

1-[2-methyl-4-(3-phenyl-2-propen-1-yl)-1-piperazinyl]-1-butanone (other name: 2-methyl AP-237);

1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP);

1-(4-cinnamyl-2,6-dimethylpiperazin-1-yl)propan-1-one (other name: AP-238);

1-methoxy-3-[4-(2-methoxy-2-phenylethyl)piperazin-1-yl]-1-phenylpropan-2-ol (other name: zipeprol);

1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP);

2-[2-[(4-ethoxyphenyl)methyl]-5-nitrobenzimidazol-1-yl]-N-ethylethanamine (other name: N-desethyl etonitazene);

2-(4-ethoxybenzyl)-5-nitro-1-(2-(piperidin-1-yl)ethyl)-1H-benzimidazole (other names: N-piperidinyl etonitazene; etonitazepipne);

2-(4-ethoxybenzyl)-5-nitro-1-[2-(pyrrolidin-1-yl)ethyl]-1H-benzimidazole (other names: N-pyrrolidino etonitazene, etonitazepyne);

2-[(4-methoxyphenyl)methyl]-5-nitro-1-(2-pyrrolidin-1-ylethyl)benzimidazole (other names: metonitazepyne, N-pyrrolidino metonitazene);

2-[(4-methoxyphenyl)methyl]-N,N-diethyl-5-nitro-1H-benzimidazole-1-ethanamine (other name:

- 31 Metonitazene);
- 32 2-methoxy-N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-acetamide (other name: Methoxyacetyl
33 fentanyl);
- 34 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzamide (other name: U-47700);
- 35 3,4-dichloro-N-{[1-(dimethylamino)cyclohexyl]methyl}benzamide (other name: AH-7921);
- 36 *3-furanyl fentanyl (other name: N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-3-carboxamide);*
- 37 *2',5'-dimethoxyfentanyl (other name: N-(1-(2,5-dimethoxyphenethyl)piperidin-4-yl)-N-*
38 *phenylpropionamide);*
- 39 7-[(3-chloro-6-methyl-5,5-dioxo-11H-benzo[c][2,1]benzothiazepin-11-yl)amino]heptanoic acid (other
40 name: Tianeptine);
- 41 Acetyl fentanyl (other name: desmethyl fentanyl);
- 42 Acetylmethadol;
- 43 Allylprodine;
- 44 Alphacetylmethadol (except levo-alphacetylmethadol, also known as levo-alpha-acetylmethadol,
45 levomethadyl acetate, or LAAM);
- 46 Alphameprodine;
- 47 Alphamethadol;
- 48 Benzethidine;
- 49 Betacetylmethadol;
- 50 Betameprodine;
- 51 Betamethadol;
- 52 Betaprodine;
- 53 Clonitazene;
- 54 Dextromoramide;
- 55 Diampromide;
- 56 Diethylthiambutene;
- 57 Difenoxin;
- 58 Dimenoxadol;
- 59 Dimepheptanol;

- 60 Dimethylthiambutene;
- 61 Dioxaphetylbutyrate;
- 62 Dipipanone;
- 63 Ethyl N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]carbamate (other name: fentanyl carbamate);
- 64 Ethylmethylthiambutene;
- 65 Etonitazene;
- 66 Etoperidone;
- 67 Furethidine;
- 68 Hydroxypethidine;
- 69 *Isovaleryl fentanyl (other name: 3-methyl-N-(1-phenethylpiperidin-4-yl)-N-phenylbutanamide);*
- 70 Ketobemidone;
- 71 Levomoramide;
- 72 Levophenacetylmorphan;
- 73 *Meta-fluorofentanyl (other name: N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)propionamide);*
- 74 *Meta-fluoroisobutyryl fentanyl (other name: N-(3-fluorophenyl)-N-(1-phenethylpiperidin-4-*
- 75 *yl)isobutyramide);*
- 76 Morpheridine;
- 77 MT-45 (1-cyclohexyl-4-(1,2-diphenylethyl)piperazine);
- 78 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide (other name: Cyclopropyl fentanyl);
- 79 N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide (other name: Tetrahydrofuranyl
- 80 fentanyl);
- 81 N-[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]-N-phenylpropanamide (other name: alpha-
- 82 methylthiofentanyl);
- 83 N-[1-(1-methyl-2-phenylethyl)-4-piperidyl]-N-phenylacetamide (other name: acetyl-alpha-
- 84 methylfentanyl);
- 85 N-{1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyl}-N-phenylpropanamide (other name: beta-
- 86 hydroxythiofentanyl);
- 87 N-[1-(2-hydroxy-2-phenyl)ethyl-4-piperidyl]-N-phenylpropanamide (other name: beta-hydroxyfentanyl);
- 88 N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]propionanilide (other names: 1-(1-methyl-2-

- 89 phenylethyl)-4-(N-propanilido) piperidine, alpha-methylfentanyl);
- 90 N-(2-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]prop-2-enamide (other name: ortho-fluoroacryl
91 fentanyl);
- 92 N-(2-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other names: 2-fluorofentanyl,
93 ortho-fluorofentanyl);
- 94 N-(2-fluorophenyl)-2-methyl-N-[1-(2-phenylethyl)piperidin-4-yl]propanamide (other name: ortho-
95 fluoroisobutyryl fentanyl);
- 96 N-(2-fluorophenyl)-N-[1-[2-(2-fluorophenyl)ethyl]piperidin-4-yl]propanamide (other names: 2'-fluoro
97 ortho-fluorofentanyl; 2'-fluoro 2-fluorofentanyl);
- 98 N-[1-[2-(4-methylphenyl)ethyl]piperidin-4-yl]-N-phenylacetamide (other name: 4'-methyl acetyl
99 fentanyl);
- 100 N-ethyl-2-[5-nitro-2-[(4-propan-2-yloxyphenyl)methyl]benzimidazol-1-yl]ethanamine (other name: N-
101 desethyl Isotonitazene);
- 102 N,3-diphenyl-N-[1-(2-phenylethyl)piperidin-4-yl]propanamide (other names: beta'-phenyl fentanyl; beta'-
103 phenyl fentanyl; 3-phenylpropanoyl fentanyl);
- 104 N-phenyl-N-[1-(2-phenylpropyl)piperidin-4-yl]propanamide (other name: beta-methyl fentanyl);
- 105 N-(2-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]butanamide (other name: ortho-fluorobutyryl
106 fentanyl; 2-fluorobutyryl fentanyl);
- 107 N-(2-methylphenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide (other names: ortho-methyl
108 acetylfentanyl; 2-methyl acetylfentanyl);
- 109 *N-(2-methylphenyl)-N-[1-(2-phenethyl)piperidin-4-yl]propanamide (other name: ortho-methylfentanyl);*
- 110 N-(4-methylphenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]propanamide (other names: para-methylfentanyl;
111 4-methylfentanyl);
- 112 N-(4-chlorophenyl)-2-methyl-N-[1-(2-phenylethyl)piperidin-4-yl]propanamide (other name: para-
113 chloroisobutyryl fentanyl);
- 114 N-(3-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: 3-fluorofentanyl);
- 115 N-[3-methyl-1-(2-hydroxy-2-phenylethyl)4-piperidyl]-N-phenylpropanamide (other name: beta-hydroxy-
116 3-methylfentanyl);
- 117 N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N-phenylpropanamide (other name: 3-methylfentanyl);

- 118 N-[3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide (other name: 3-methylthiofentanyl);
- 119 N-(4-chlorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other names: para-chlorofentanyl,
- 120 4-chlorofentanyl);
- 121 N-(4-fluorophenyl)-2-methyl-N-[1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: para-
- 122 fluoroisobutyryl fentanyl);
- 123 N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]furan-2-carboxamide (other name: para-fluoro
- 124 furanyl fentanyl);
- 125 N-(4-fluorophenyl)-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: para-
- 126 fluorobutyrylfentanyl);
- 127 N-(4-fluorophenyl)-N-1-(2-phenylethyl)-4-piperidinyl]-propanamide (other name: para-fluorofentanyl);
- 128 N,N-diethyl-2-[2-[(4-fluorophenyl)methyl]-5-nitrobenzimidazol-1-yl] ethanamine (other name:
- 129 Flunitazene);
- 130 N,N-diethyl-2-(2-(4-isopropoxybenzyl)-5-nitro-1H-benzimidazol-1-yl)ethan-1-amine (other name:
- 131 Isotonitazene);
- 132 N,N-diethyl-2-[[4-ethoxyphenyl] methyl]-1H-benzimidazol-1-yl}-ethan-1-amine (other names: Etazene,
- 133 Desnitroetonitazene);
- 134 N,N-diethyl-2-[(4-methoxyphenyl)methyl]-1H-benzimidazole-1-ethanamine (other name:
- 135 Metodesnitazene);
- 136 N,N-diethyl-2-[5-nitro-2-(4-propoxybenzyl)-1H-benzimidazol-1-yl]ethanamine (other name:
- 137 Protonitazene);
- 138 N-phenyl-N-[1-(2-phenylmethyl)-4-piperidinyl]-2-furancarboxamide (other name: N-benzyl Furanyl
- 139 norfentanyl);
- 140 N-phenyl-N-(4-piperidinyl)-propanamide (other name: Norfentanyl);
- 141 Noracymethadol;
- 142 Norlevorphanol;
- 143 Normethadone;
- 144 Norpipanone;
- 145 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-furancarboxamide (other name: Furanyl fentanyl);
- 146 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-2-propenamide (other name: Acryl fentanyl);

- 147 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-butanamide (other name: butyryl fentanyl);
- 148 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-pentanamide (other name: Pentanoyl fentanyl);
- 149 N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide (other name: thiofentanyl);
- 150 N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]thiophene-2-carboxamide (other name: thiophene fentanyl);
- 151 *Ortho-fluorofuranyl fentanyl (other name: N-(2-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-*
- 152 *carboxamide);*
- 153 *Para-methoxyfuranyl fentanyl (other name: N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)furan-2-*
- 154 *carboxamide);*
- 155 *Para-methylcyclopropyl fentanyl (other name: N-(4-methylphenyl)-N-(1-phenethylpiperidin-4-*
- 156 *yl)cyclopropanecarboxamide);*
- 157 Phenadoxone;
- 158 Phenampromide;
- 159 Phenomorphan;
- 160 Phenoperidine;
- 161 Piritramide;
- 162 Proheptazine;
- 163 Properidine;
- 164 Propiram;
- 165 Racemoramide;
- 166 Tilidine;
- 167 Trimeperidine;
- 168 N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]-1,3-benzodioxole-5-carboxamide (other name:
- 169 Benzodioxole fentanyl);
- 170 3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methylbenzamide (other name: U-49900);
- 171 2-methoxy-N-(2-methylphenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide (other names: ortho-methyl
- 172 methoxyacetylfentanyl; 2-methyl methoxyacetyl fentanyl);
- 173 2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]butanamide (other names: 2-methyl butyryl
- 174 fentanyl, alpha'-methyl butyryl fentanyl);
- 175 2-[2-[(4-butoxyphenyl)methyl]-5-nitrobenzimidazol-1-yl]-N,N-diethylethanamine (other name:

- 176 Butonitazene);
- 177 2-(2,4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methyl acetamide (other name: U-48800);
- 178 2-(3,4-dichlorophenyl)-N-[2-(dimethylamino)cyclohexyl]-N-methyl acetamide (other name: U-51754);
- 179 2-(4-isopropoxybenzyl)-5-nitro-1-[2-(pyrrolidin-1-yl)ethyl]-1H-benzo[d]imidazole (other name: N-
- 180 Pyrrolidino Isotonitazene);
- 181 N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)-4-piperidiny]-acetamide (other name: Ocfentanil);
- 182 N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)pentanamide (other names: para-fluoro valeryl fentanyl,
- 183 para-fluoro pentanoyl fentanyl);
- 184 N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]acetamide (other name: para-fluoroacetyl
- 185 fentanyl);
- 186 N-(4-methoxyphenyl)-N-[1-(2-phenylethyl)-4-piperidiny]-butanamide (other name: 4-
- 187 methoxybutyrylfentanyl);
- 188 N-phenyl-2-methyl-N-[1-(2-phenylethyl)-4-piperidiny]-propanamide (other name: Isobutyryl fentanyl);
- 189 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-cyclopentanecarboxamide (other name: Cyclopentyl
- 190 fentanyl);
- 191 N-phenyl-N-(1-methyl-4-piperidiny)-propanamide (other name: N-methyl norfentanyl);
- 192 N-phenyl-N-(1-propionyl-4-piperidiny)-propanamide (other name: N-propionyl norfentanyl);
- 193 N-[2-(dimethylamino)cyclohexyl]-N-methyl-1,3-benzodioxole-5-carboxamide (other names: 3,4-
- 194 methylenedioxy U-47700 or 3,4-MDO-U-47700);
- 195 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-2-butenamide (other name: Crotonyl fentanyl);
- 196 N-phenyl-N-[4-phenyl-1-(2-phenylethyl)-4-piperidiny]-propanamide (other name: 4-phenylfentanyl);
- 197 N-phenyl-N-[1-(2-phenylethyl)-4-piperidiny]-benzamide (other names: Phenyl fentanyl, Benzoyl
- 198 fentanyl);
- 199 N-[2-(dimethylamino)cyclohexyl]-N-phenylfuran-2-carboxamide (other name: Furanyl UF-17);
- 200 N-[2-(dimethylamino)cyclohexyl]-N-phenylpropionamide (other name: UF-17);
- 201 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-isopropyl-benzamide (other name: Isopropyl U-47700);
- 202 5-nitro-2-(4-propoxybenzyl)-1-[2-(pyrrolidin-1-yl)ethyl]-1H-benzo[d]imidazole (other names: N-
- 203 Pyrrolidino Protonitazene, Protonitazepyne).
- 204 2. Any of the following opium derivatives, their salts, isomers and salts of isomers, unless specifically

205 excepted, whenever the existence of these salts, isomers and salts of isomers is possible within the specific
206 chemical designation:

207 Acetorphine;

208 Acetyldihydrocodeine;

209 Benzylmorphine;

210 Codeine methylbromide;

211 Codeine-N-Oxide;

212 Cyprenorphine;

213 Desomorphine;

214 Dihydromorphine;

215 Drotebanol;

216 Etorphine;

217 Heroin;

218 Hydromorphenol;

219 Methyldesorphine;

220 Methyldihydromorphine;

221 Morphine methylbromide;

222 Morphine methylsulfonate;

223 Morphine-N-Oxide;

224 Myorphine;

225 Nicocodeine;

226 Nicomorphine;

227 Normorphine;

228 Pholcodine;

229 Thebacon.

230 3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture, or
231 preparation, which contains any quantity of the following hallucinogenic substances, or which contains any
232 of its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers, and salts of isomers
233 is possible within the specific chemical designation (for purposes of this subdivision only, the term "isomer"

234 includes the optical, position, and geometric isomers):
235 3-[2-(diethylamino)ethyl]-1H-indol-4-ol (other names: 4-hydroxy-N,N-diethyltryptamine; 4-hydroxy
236 DET; ethocin);
237 [3-[2-(diethylamino)ethyl]-1H-indol-4-yl] acetate (other names: 4-acetoxy-N,N-diethyltryptamine; 4-
238 acetoxy DET; 4-AcO-DET; ethacetin);
239 3-methylmethcathinone (other names: 3-MMC; metaphedrone; 2-(methylamino)-1-(3-
240 methylphenyl)propan-1-one);
241 Alpha-ethyltryptamine (some trade or other names: Monase; a-ethyl-1H-indole-3-ethanamine; 3-2-
242 aminobutyl] indole; a-ET; AET);
243 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names: 2-4-bromo-2,5-dimethoxyphenyl]-1-
244 aminoethane;alpha-desmethyl DOB; 2C-B; Nexus);
245 3,4-methylenedioxy amphetamine;
246 5-methoxy-3,4-methylenedioxy amphetamine;
247 3,4,5-trimethoxy amphetamine;
248 Alpha-methyltryptamine (other name: AMT);
249 Bufotenine;
250 Diethyltryptamine;
251 Dimethyltryptamine;
252 4-methyl-2,5-dimethoxyamphetamine;
253 2,5-dimethoxy-4-ethylamphetamine (DOET);
254 4-fluoro-N-ethylamphetamine;
255 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);
256 Ibogaine;
257 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);
258 Lysergic acid diethylamide;
259 Mescaline;
260 Parahexyl (some trade or other names: 3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-
261 dibenzo [b,d] pyran; Synhexyl);
262 Peyote;

- 263 N-ethyl-3-piperidyl benzilate;
- 264 N-methyl-3-piperidyl benzilate;
- 265 *N,N-dipropyl -1H-indole-3-ethanamine (other names: Dipropyltryptamine; N,N-DPT);*
- 266 Psilocybin;
- 267 Psilocyn;
- 268 Salvinorin A;
- 269 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy- α -methylphenethylamine; 2,5-
- 270 DMA);
- 271 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts and
- 272 salts of isomers;
- 273 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl- α -methyl-3,4
- 274 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);
- 275 N-hydroxy-3,4-methylenedioxyamphetamine (some other names: N-hydroxy- α -methyl-
- 276 3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);
- 277 4-bromo-2,5-dimethoxyamphetamine (some trade or other names: 4-bromo-2,5-dimethoxy- α -
- 278 methylphenethylamine; 4-bromo-2,5-DMA);
- 279 4-methoxyamphetamine (some trade or other names: 4-methoxy- α -methylphenethylamine;
- 280 paramethoxyamphetamine; PMA);
- 281 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine, (1-
- 282 phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);
- 283 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP);
- 284 Thiophene analog of phencyclidine (some other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienyl
- 285 analog of phencyclidine, TPCP, TCP);
- 286 1-1-(2-thienyl)cyclohexylpyrrolidine (other name: TCPy);
- 287 3,4-methylenedioxypyrovalerone (other name: MDPV);
- 288 4-methylmethcathinone (other names: mephedrone, 4-MMC);
- 289 3,4-methylenedioxymethcathinone (other name: methylone);
- 290 Naphthylpyrovalerone (other name: naphyrone);
- 291 4-fluoromethcathinone (other names: flephedrone, 4-FMC);

- 292 4-methoxymethcathinone (other names: methedrone; bk-PMMA);
- 293 Ethcathinone (other name: N-ethylcathinone);
- 294 3,4-methylenedioxyethcathinone (other name: ethylone);
- 295 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone);
- 296 N,N-dimethylcathinone (other name: metamfepramone);
- 297 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);
- 298 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);
- 299 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);
- 300 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);
- 301 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);
- 302 3-fluoromethcathinone (other name: 3-FMC);
- 303 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);
- 304 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);
- 305 4-Methylethcathinone (other name: 4-MEC);
- 306 4-Ethylmethcathinone (other name: 4-EMC);
- 307 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);
- 308 Beta-keto-methylbenzodioxolylpentanamine (other names: Pentylone, bk-MBDP);
- 309 Alpha-methylamino-butyrophenone (other name: Buphedrone);
- 310 Alpha-methylamino-valerophenone (other name: Pentedrone);
- 311 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);
- 312 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);
- 313 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I, 25I-
314 NBOMe, 2C-I-NBOMe);
- 315 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);
- 316 4-Fluoromethamphetamine (other name: 4-FMA);
- 317 4-Fluoroamphetamine (other name: 4-FA);
- 318 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);
- 319 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);
- 320 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);

- 321 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);
- 322 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);
- 323 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);
- 324 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);
- 325 (2-aminopropyl)benzofuran (other name: APB);
- 326 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);
- 327 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 2C-C-NBOMe,
- 328 25C-NBOMe, 25C);
- 329 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 2C-B-NBOMe,
- 330 25B-NBOMe, 25B);
- 331 Acetoxymethyltryptamine (other names: ~~AcO-Psilocin~~, AcO-DMT, Psilacetin);
- 332 Benocyclidine (other names: BCP, BTCP);
- 333 Alpha-pyrrolidinobutiophenone (other name: alpha-PBP);
- 334 3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
- 335 4-bromomethcathinone (other name: 4-BMC);
- 336 4-chloromethcathinone (other name: 4-CMC);
- 337 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
- 338 Alpha-Pyrrolidinohexiophenone (other name: alpha-PHP);
- 339 Alpha-Pyrrolidinoheptiophenone (other name: PV8);
- 340 5-methoxy-N,N-methylisopropyltryptamine (other name: 5-MeO-MIPT);
- 341 Beta-keto-N,N-dimethylbenzodioxolylbutanamine (other names: Dibutylone, bk-DMBDB);
- 342 Beta-keto-4-bromo-2,5-dimethoxyphenethylamine (other name: bk-2C-B);
- 343 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-pentanone (other name: N-ethylpentylone);
- 344 1-[1-(3-methoxyphenyl)cyclohexyl]piperidine (other name: 3-methoxy PCP);
- 345 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine (other name: 4-methoxy PCP);
- 346 4-Chloroethcathinone (other name: 4-CEC);
- 347 3-Methoxy-2-(methylamino)-1-(4-methylphenyl)-1-propanone (other name: Mexedrone);
- 348 1-propionyl lysergic acid diethylamide (other name: 1P-LSD);
- 349 (2-Methylaminopropyl)benzofuran (other name: MAPB);

- 350 1-(1,3-benzodioxol-5-yl)-2-(dimethylamino)-1-pentanone (other names: N,N-Dimethylpentylone,
351 Dipentylone);
- 352 1-(4-methoxyphenyl)-2-(pyrrolidin-1-yl)octan-1-one (other name: 4-methoxy-PV9);
- 353 3,4-tetramethylene-alpha-pyrrolidinovalerophenone (other name: TH-PVP);
- 354 4-allyloxy-3,5-dimethoxyphenethylamine (other name: Allylescaline);
- 355 4-Bromo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25B-NBOH);
- 356 4-chloro-alpha-methylamino-valerophenone (other name: 4-chloropentedrone);
- 357 4-chloro-alpha-Pyrrolidinovalerophenone (other name: 4-chloro-alpha-PVP);
- 358 4-fluoro-alpha-Pyrrolidinoheptiophenone (other name: 4-fluoro-PV8);
- 359 4-hydroxy-N,N-diisopropyltryptamine (other name: 4-OH-DIPT);
- 360 4-methyl-alpha-ethylaminopentiophenone;
- 361 4-methyl-alpha-Pyrrolidinohexiophenone (other name: MPHP);
- 362 5-methoxy-N,N-dimethyltryptamine (other name: 5-MeO-DMT);
- 363 5-methoxy-N-ethyl-N-isopropyltryptamine (other name: 5-MeO-EIPT);
- 364 6-ethyl-6-nor-lysergic acid diethylamide (other name: ETH-LAD);
- 365 6-allyl-6-nor-lysergic acid diethylamide (other name: AL-LAD);
- 366 (N-methyl aminopropyl)-2,3-dihydrobenzofuran (other name: MAPDB);
- 367 2-(methylamino)-2-phenyl-cyclohexanone (other name: Deschloroketamine);
- 368 2-(ethylamino)-2-phenyl-cyclohexanone (other name: deschloro-N-ethyl-ketamine);
- 369 2-methyl-1-(4-(methylthio)phenyl)-2-morpholinopropiophenone (other name: MMMP);
- 370 Alpha-ethylaminohexanophenone (other name: N-ethylhexedrone);
- 371 N-ethyl-1-(3-methoxyphenyl)cyclohexylamine (other name: 3-methoxy-PCE);
- 372 4-fluoro-alpha-pyrrolidinohexiophenone (other name: 4-fluoro-alpha-PHP);
- 373 N-ethyl-1,2-diphenylethylamine (other name: Ephenedine);
- 374 2,5-dimethoxy-4-chloroamphetamine (other name: DOC);
- 375 3,4-methylenedioxy-N-tert-butylcathinone;
- 376 Alpha-pyrrolidinoisohexiophenone (other name: alpha-PiHP);
- 377 1-[1-(3-hydroxyphenyl)cyclohexyl]piperidine (other name: 3-hydroxy PCP);
- 378 4-acetyloxy-N,N-diallyltryptamine (other name: 4-AcO-DALT);

- 379 4-hydroxy-N,N-methylisopropyltryptamine (other name: 4-hydroxy-MiPT);
- 380 3,4-Methylenedioxy-alpha-pyrrolidinohexanophenone (other name: MDPHP);
- 381 5-methoxy-N,N-dibutyltryptamine (other name: 5-methoxy-DBT);
- 382 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-1-butanone (other names: Eutylone, bk-EBDB);
- 383 1-(1,3-benzodioxol-5-yl)-2-(butylamino)-1-pentanone (other name: N-butylpentylone);
- 384 N-benzyl-3,4-dimethoxyamphetamine (other name: N-benzyl-3,4-DMA);
- 385 1-(benzo[d][1,3]dioxol-5-yl)-2-(sec-butylamino)pentan-1-one (other name: N-sec-butyl Pentylone);
- 386 1-cyclopropionyl lysergic acid diethylamide (other name: 1cP-LSD);
- 387 2-(ethylamino)-1-phenylheptan-1-one (other name: N-ethylheptedrone);
- 388 (2-ethylaminopropyl)benzofuran (other name: EAPB);
- 389 4-ethyl-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25E-NBOH);
- 390 2-fluoro-Deschloroketamine (other name: 2-(2-fluorophenyl)-2-(methylamino)-cyclohexanone);
- 391 4-hydroxy-N-ethyl-N-propyltryptamine (other name: 4-hydroxy-EPT);
- 392 2-(isobutylamino)-1-phenylhexan-1-one (other names: N-Isobutyl Hexedrone, alpha-
- 393 isobutylaminohexanphenone);
- 394 1-(4-methoxyphenyl)-N-methylpropan-2-amine (other names: para-Methoxymethamphetamine, PMMA);
- 395 N-ethyl-1-(3-hydroxyphenyl)cyclohexylamine (other name: 3-hydroxy-PCE);
- 396 N-heptyl-3,4-dimethoxyamphetamine (other name: N-heptyl-3,4-DMA);
- 397 N-hexyl-3,4-dimethoxyamphetamine (other name: N-hexyl-3,4-DMA);
- 398 4-fluoro-3-methyl-alpha-pyrrolidinovalerophenone (other name: 4-fluoro-3-methyl-alpha-PVP);
- 399 4-fluoro-alpha-methylamino-valerophenone (other name: 4-fluoropentedrone);
- 400 N-(1,4-dimethylpentyl)-3,4-dimethoxyamphetamine (other name: N-(1,4-dimethylpentyl)-3,4-DMA);
- 401 4,5-methylenedioxy-N,N-diisopropyltryptamine (other name: 4,5-MDO-DiPT);
- 402 Alpha-pyrrolidinocyclohexanophenone (other name: alpha-PCYP);
- 403 3,4-methylenedioxy-alpha-pyrrolidinoheptiophenone (other name: MDPV8);
- 404 4-chloro-alpha-methylaminobutiophenone (other name: 4-chloro Buphedrone);
- 405 1-(1,3-benzodioxol-5-yl)-2-(cyclohexylamino)butan-1-one (other names: Cybutylone, N-cyclohexyl
- 406 Butylone);
- 407 1-(1,3-benzodioxol-5-yl)-2-(propylamino)-1-butanone (other names: 3,4-Methylenedioxy-alpha-

- 408 propylaminobutiophenone; N-propyl butylone);
- 409 1-[1-(3-chlorophenyl)cyclohexyl]-piperidine (other names: 3-Chloro Phencyclidine, 3Cl-PCP, 3-chloro
- 410 PCP);
- 411 1-[1-(3-fluorophenyl)cyclohexyl]piperidine (other names: 3-fluoro Phencyclidine, 3F-PCP);
- 412 1-(7-methoxy-1,3-benzodioxol-5-yl)propan-2-amine (other names: 5-methoxy-3,4-
- 413 methylenedioxyamphetamine, 3-methoxy MDA, MMDA);
- 414 2-(ethylamino)-1-phenylpentan-1-one (other names: N-ethylpentedrone, alpha-ethylaminopentiophenone);
- 415 2-(ethylamino)-2-(2-fluorophenyl)-cyclohexanone (other names: 2-fluoro-2-oxo PCE, 2-fluoro
- 416 NENDCK);
- 417 3,4-methylenedioxy-alpha-cyclohexylaminopropiophenone (other name: Cyputylone);
- 418 3,4-methylenedioxy-alpha-cyclohexylmethylaminopropiophenone (other name: 3,4-Methylenedioxy-N,N-
- 419 cyclohexylmethcathinone);
- 420 3,4-methylenedioxy-alpha-isopropylaminobutiophenone (other name: N-isopropyl butylone);
- 421 4-chloro-N-butylcathinone (other names: 4-chlorobutylcathinone, para-chloro-N-butylcathinone);
- 422 4-hydroxy-N-methyl-N-ethyltryptamine (other names: 4-hydroxy MET, Metocin);
- 423 4-methallyloxy-3,5-dimethoxyphenethylamine (other name: Methallylescaline);
- 424 Alpha-pyrrolidino-2-phenylacetophenone (other name: alpha-D2PV);
- 425 1-(3,5-Dimethoxy-4-propoxyphenyl)-2-propanamine (other names: 4-propoxy-3,5-DMA, 3C-P, 1-(3,5-
- 426 Dimethoxy-4-propoxyphenyl)propan-2-amine);
- 427 2-(5-methoxy-1H-indol-3-yl)ethanamine (other names: 5-methoxytryptamine, 5-MeOT);
- 428 1-(1,3-benzodioxol-5-yl)-2-(isobutylamino)-1-pentanone (other name: N-isobutylpentylone);
- 429 1-(1,3-benzodioxyl-5-yl)-2-(tert-butylamino)-1-pentanone (other name: N-tert-butyl pentylone);
- 430 1-Phenyl-N-propylcyclohexanamine (other names: N-(1-phenylcyclohexyl)propanamine, PCPr);
- 431 1-[(4-fluorophenyl)methyl]-4-methylpiperazine (other names: 4-fluoro-MBZP, 4-fluoro
- 432 methylbenzylpiperazine);
- 433 4-fluoro-alpha-pyrrolidinoisohexiophenone (other name: 4-fluoro-alpha-PiHP);
- 434 8-bromo-1-methyl-6-pyridin-2-yl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine (other name: pyrazolam).
- 435 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture or
- 436 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:

5-(2-chlorophenyl)-1,3-dihydro-3-methyl-7-nitro-2H-1,4-benzodiazepin-2-one (other name: Meclonazepam);

7-Bromo-5-(2-chlorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one (other name: phenazepam);

7-chloro-5-(2-fluorophenyl)-1,3-dihydro-1,4-benzodiazepin-2-one (other name: Norfludiazepam);

Bromazolam;

Clonazolam;

Deschloroetizolam;

Etizolam;

Flualprazolam;

Flubromazepam;

Flubromazolam;

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

Mecloqualone;

Methaqualone;

6-(4-chlorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine (other names: 4'-chloro Deschloroalprazolam, 4'Cl-Deschloroalprazolam);

7-bromo-5-phenyl-1,3-dihydro-1,4-benzodiazepin-2-one (other names: Desalkylgidazepam, Bromonordiazepam);

7-chloro-5-(2-chlorophenyl)-1-methyl-3H-1,4-benzodiazepin-2-one (other names: Diclazepam, 2-Chlorodiazepam);

8-bromo-6-(2-chlorophenyl)-1-methyl-4H-[1,2,4]triazolo[4,3-a][1,4]benzodiazepine (other names: Clobromazolam, Phenazolam).

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);

- 466 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,5-dihydro-5-
467 phenyl-2-oxazolamine);
- 468 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone, 2-
469 aminopropiophenone, norephedrone), and any plant material from which Cathinone may be derived;
- 470 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
- 471 Ethylamphetamine;
- 472 Ethyl phenyl(piperidin-2-yl)acetate (other name: Ethylphenidate);
- 473 Fenethylamine;
- 474 Methcathinone (some other names: 2-(methylamino)-propionophenone; alpha-(methylamino)-
475 propionophenone; 2-(methylamino)-1-phenylpropan-1-one; alpha-N-methylaminopropionophenone;
476 monomethylpropion; ephedrone; N-methylcathinone; methylcathinone; AL-464; AL-422; AL-463 and UR
477 1432);
- 478 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
- 479 N-methyl-1-(thiophen-2-yl)propan-2-amine (other name: methiopropamine);
- 480 N,N-dimethylamphetamine (other names: N,N-alpha-trimethyl-benzeneethanamine, N,N-alpha-
481 trimethylphenethylamine);
- 482 N-phenyl-N'-(3-(1-phenylpropan-2-yl)-1,2,3-oxadiazol-3-ium-5-yl)carbamimidate (other name:
483 mesocarb);
- 484 Methyl 2-(4-fluorophenyl)-2-(2-piperidinyl)acetate (other name: 4-fluoromethylphenidate);
- 485 Isopropyl-2-phenyl-2-(2-piperidinyl)acetate (other name: Isopropylphenidate);
- 486 2-(3-chlorophenyl)-3-methylmorpholine (other name: 3-chlorophenmetrazine);
- 487 4-chloro-N,N-dimethylcathinone;
- 488 3,4-methylenedioxy-N-benzylcathinone (other name: BMDP);
- 489 4-methylmethamphetamine (other names: N-alpha,4-trimethyl-benzeneethanamine, 4-MMA);
- 490 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine (other names: 4,4'-Dimethylaminorex, 4,4'-
491 DMAR);
- 492 7-[(10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-yl)amino]heptanoic acid (other name: amineptine).
- 493 6. Any substance that contains one or more cannabimimetic agents or that contains their salts, isomers,
494 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 classes:

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 any extent;

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 any extent;

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 extent;

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 extent;

N-(adamantyl)-indole-3-carboxamide 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 extent; and

N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring, whether or not further substituted on the indazole ring to any extent, whether or not substituted on the adamantyl ring to any extent.

b. The term "cannabimimetic agents" includes:

Methyl N-[(5-methyl-1H-indazol-3-yl)carbonyl]-3-methyl-valinate (other name: MDMB-5Me-INACA);

N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indazole-3-carboxamide (other name: ADB-INACA);

N-cyclohexyl-2-(1-pentylindol-3-yl)acetamide (other names: cyclohexyl-PIATA, CH-PIACA, CH-PIATA);

- 524 *N*-(1-amino-3-methyl-1-oxobutan-2-yl)-3-(dimethylsulfamoyl)-4-methylbenzamide (other name: AB-
525 MDMSBA);
- 526 Methyl 2-[[1-(4-fluorobutyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate (other names: 4F-MDMB-
527 BUTICA; 4F-MDMB-BICA);
- 528 5-Pentyl-2-(2-phenylpropan-2-yl)pyrido[4,3-b]indol-1-one (other names: CUMYL-PEGACLONE; SGT-
529 151);
- 530 Ethyl 2-[[1-(5-fluoropentyl)indole-3-carbonyl]amino]-3,3-dimethyl-butanoate (other names: 5F-EDMB-
531 PICA; 5F-EDMB-2201);
- 532 5-bromo-N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-indazole-3-carboxamide (other name: ADB-5Br-
533 INACA);
- 534 5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);
- 535 5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);
- 536 5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);
- 537 5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);
- 538 1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);
- 539 1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);
- 540 1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);
- 541 1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);
- 542 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);
- 543 (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tet
544 rahydrobenzo[c]chromen-1-ol (other name: HU-210);
- 545 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);
- 546 1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);
- 547 1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);
- 548 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);
- 549 1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);
- 550 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);
- 551 1-((N-methylpiperidin-2-yl)methyl)-3-(1-naphthoyl)indole (other name: AM-1220);
- 552 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);

- 553 1-[(N-methylpiperidin-2-yl)methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);
- 554 Pravadoline (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone (other name:
- 555 WIN 48,098);
- 556 1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);
- 557 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);
- 558 1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);
- 559 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other names: XLR-11, 5-fluoro-
- 560 UR-144);
- 561 N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);
- 562 N-adamantyl-1-pentylindazole-3-carboxamide (other names: AKB48, APINACA);
- 563 1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);
- 564 (8-quinoliny)(1-pentylindol-3-yl)carboxylate (other name: PB-22);
- 565 (8-quinoliny)(1-(5-fluoropentyl)indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);
- 566 (8-quinoliny)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);
- 567 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-5-bromo-1-butyldiazole-3-carboxamide (other name: ADB-
- 568 5'Br-BUTINACA);
- 569 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);
- 570 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name: AB-
- 571 FUBINACA);
- 572 1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);
- 573 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: ADB-
- 574 PINACA);
- 575 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name: AB-
- 576 CHMINACA);
- 577 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name: 5-fluoro-
- 578 AB-PINACA);
- 579 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other names:
- 580 ADB-CHMINACA, MAB-CHMINACA);
- 581 Methyl N-[(5-bromo-1H-indazol-3-yl)carbonyl]-3-methyl-valinate (other name: MDMB-5Br-INACA);

582 Methyl-2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name: 5-fluoro-
583 AMB);

584 1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);

585 1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);

586 1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name MAM-2201);

587 N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamide
588 (other name: ADB-FUBINACA);

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

591 Methyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names: 5-fluoro-
592 ADB, 5-Fluoro-MDMB-PINACA);

593 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other names:
594 AMB-FUBINACA, FUB-AMB);

595 N-(adamantan-1-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (other names: FUB-AKB48, 5F-
596 APINACA);

597 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (other name: 5F-AKB48);

598 N-(adamantan-1-yl)-1-(5-chloropentyl) indazole-3-carboxamide (other name: 5-chloro-AKB48);

599 Naphthalen-1-yl 1-pentyl-1H-indazole-3-carboxylate (other name: SDB-005);

600 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indole-3-carboxamide (other name: AB-
601 CHMICA);

602 1-pentyl-N-(phenylmethyl)-1H-indole-3-carboxamide (other name: SDB-006);

603 Quinolin-8-yl 1-(4-fluorobenzyl)-1H-indole-3-carboxylate (other name: FUB-PB-22);

604 Methyl N-[1-(cyclohexylmethyl)-1H-indole-3-carbonyl]valinate (other name: MMB-CHMICA);

605 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name: 5-
606 fluoro-ADB-PINACA);

607 1-(4-cyanobutyl)-N-(1-methyl-1-phenylethyl)-1H-indazole-3-carboxamide (other name: 4-cyano
608 CUMYL-BUTINACA);

609 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other names: 5-fluoro
610 MDMB-PICA, 5F-MDMB-PICA);

- 611 Ethyl 2-({1-[(4-fluorophenyl)methyl]-1H-indazole-3-carbonyl}amino)-3-methylbutanoate (other name:
612 EMB-FUBINACA);
- 613 Ethyl-3,3-dimethyl-2-[(1-(pent-4-enylindazole-3-carbonyl)amino]butanoate (other name: EDMB-4en-
614 PINACA);
- 615 Methyl 2-[1-4-fluorobutyl]-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name: 4-fluoro-
616 MDMB-BUTINACA);
- 617 1-(5-fluoropentyl)-N-(1-methyl-1-phenylethyl)-1H-indole-3-carboxamide (other name: 5-fluoro CUMYL-
618 PICA);
- 619 Methyl 2-[1-(pent-4-enyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name: MDMB-
620 4en-PINACA);
- 621 Methyl 2-({1-[(4-fluorophenyl)methyl]-1H-indole-3-carbonyl}amino)-3-methylbutanoate (other names:
622 MMB-FUBICA, AMB-FUBICA);
- 623 Methyl 2-[1-(4-penten-1-yl)-1H-indole-3-carboxamido]-3-methylbutanoate (other names: MMB022,
624 MMB-4en-PICA);
- 625 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-methylbutanoate (other name: MMB 2201);
- 626 Methyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-phenylpropanoate (other name: 5-fluoro-MPP-
627 PICA);
- 628 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyldiazole-3-carboxamide (other name: ADB-
629 BUTINACA);
- 630 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-chloropentyl)indazole-3-carboxamide (other name: 5-chloro-
631 AB-PINACA);
- 632 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)-1H-indazole-3-carboxamide (other names: 5F-CUMYL-
633 PINACA, 5-fluoro CUMYL-PINACA, CUMYL-5F-PINACA);
- 634 1-(5-fluoropentyl)-N-(2-phenylpropan-2-yl)pyrrolo[2,3-b]pyridine-3-carboxamide (other name: 5F-
635 CUMYL-P7AICA);
- 636 Ethyl 2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other names: 5F-
637 EDMB-PINACA, 5-fluoro EDMB-PINACA);
- 638 Ethyl-2-[1-(5-fluoropentyl)-1H-indazole-3-carboxamido]-3-methylbutanoate (other names: 5-fluoro-
639 EMB-PINACA, 5F-AEB);

- 640 Ethyl 2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3-methylbutanoate (other name: 5-fluoro-EMB-
641 PICA);
- 642 Ethyl-2-[1-(5-fluoropentyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name: 5-fluoro
643 EDMB-PICA);
- 644 Methyl 2-[1-(4-fluorobutyl)-1H-indole-3-carboxamido]-3,3-dimethylbutanoate (other name: 4-fluoro-
645 MDMB-BUTICA);
- 646 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate (other names:
647 MDMB-CHMICA, MMB-CHMINACA);
- 648 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(pent-4-enyl)indazole-3-carboxamide (other name: ADB-
649 4en-PINACA);
- 650 Ethyl 2-[1-pentyl-1H-indazole-3-carboxamido]-3,3-dimethylbutanoate (other name: EDMB-PINACA);
- 651 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-phenethyl-1H-indazole-3-carboxamide (other name: ADB-
652 PHETINACA);
- 653 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indole-3-acetamide (other names:
654 ADB-FUBIATA, AD-18, FUB-ACADB);
- 655 Methyl N-(1H-indazol-3-ylcarbonyl)-3-methyl-valinate (other name: MDMB-INACA);
- 656 Methyl-2-(1-butyl-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (other name: MDMB-
657 BUTINACA).
- 658 **2. That the provisions of this act may result in a net increase in periods of imprisonment or**
659 **commitment. Pursuant to § 30-19.1:4 of the Code of Virginia, the estimated amount of the necessary**
660 **appropriation is \$0 for periods of imprisonment in state adult correctional facilities and cannot be**
661 **determined for periods of commitment to the custody of the Department of Juvenile Justice.**