

## 1 VIRGINIA ACTS OF ASSEMBLY — CHAPTER

2 *An Act to amend and reenact § 54.1-3446 of the Code of Virginia, relating to Drug Control Act;*  
3 *Schedule I drugs; addition of substances.*

4 [S 480]

5 Approved

6 **Be it enacted by the General Assembly of Virginia:**7 **1. That § 54.1-3446 of the Code of Virginia is amended and reenacted as follows:**8 **§ 54.1-3446. Schedule I.**

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

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

13 Acetylmethadol;

14 Allylprodine;

15 Alphacetylmethadol (except levo-alphacetylmethadol, also known as levo-alpha-acetylmethadol,  
16 levomethadyl acetate, or LAAM);

17 Alphameprodine;

18 Alphamethadol;

19 Benzethidine;

20 Betacetylmethadol;

21 Betameprodine;

22 Betamethadol;

23 Betaprodine;

24 Clonitazene;

25 Dextromoramide;

26 Diampromide;

27 Diethylthiambutene;

28 Difenoxin;

29 Dimenoxadol;

30 Dimepheptanol;

31 Dimethylthiambutene;

32 Dioxaphetylbutyrate;

33 Dipipanone;

34 Ethylmethylthiambutene;

35 Etonitazene;

36 Etoxidine;

37 Furethidine;

38 Hydroxypethidine;

39 Ketobemidone;

40 Levomoramide;

41 Levophenacymorphan;

42 Morpheridine;

43 Noracymethadol;

44 Norlevorphanol;

45 Normethadone;

46 Norpipanone;

47 Phenadoxone;

48 Phenampromide;

49 Phenomorphan;

50 Phenoperidine;

51 Piritramide;

52 Proheptazine;

53 Properidine;

54 Propiram;

55 Racemoramide;

56 Tilidine;

57 Trimeperidine.

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

61 Acetorphine;  
62 Acetyldihydrocodeine;  
63 Benzylmorphine;  
64 Codeine methylbromide;  
65 Codeine-N-Oxide;  
66 Cyprenorphine;  
67 Desomorphine;  
68 Dihydromorphine;  
69 Drotebanol;  
70 Etorphine;  
71 Heroin;  
72 Hydromorphanol;  
73 Methyldesorphine;  
74 Methyldihydromorphine;  
75 Morphine methylbromide;  
76 Morphine methylsulfonate;  
77 Morphine-N-Oxide;  
78 Myrophine;  
79 Nicocodeine;  
80 Nicomorphine;  
81 Normorphine;  
82 Pholcodine;  
83 Thebacon.

84 3. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture,  
85 or preparation, which contains any quantity of the following hallucinogenic substances, or which  
86 contains any of its salts, isomers, and salts of isomers, whenever the existence of such salts, isomers,  
87 and salts of isomers is possible within the specific chemical designation (for purposes of this subdivision  
88 only, the term "isomer" includes the optical, position, and geometric isomers):

89 Alpha-ethyltryptamine (some trade or other names: Monase; a-ethyl-1H-indole-3-ethanamine;  
90 3-2-aminobutyl] indole; a-ET; AET);

91 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:  
92 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B; Nexus);

93 4-Bromo-2,5-dimethoxyphenethylamine (some trade or other names:  
94 2-4-bromo-2,5-dimethoxyphenyl]-1-aminoethane; alpha-desmethyl DOB; 2C-B; Nexus);

95 3,4-methylenedioxy amphetamine;

96 5-methoxy-3,4-methylenedioxy amphetamine;

97 3,4,5-trimethoxy amphetamine;

98 Alpha-methyltryptamine (other name: AMT);

99 Bufotenine;

100 Diethyltryptamine;

101 Dimethyltryptamine;

102 4-methyl-2,5-dimethoxyamphetamine;

103 2,5-dimethoxy-4-ethylamphetamine (DOET);

104 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

105 Ibogaine;

106 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT);

107 Lysergic acid diethylamide;

108 Mescaline;

109 Parahexyl (some trade or other names: 3-Hexyl-1-hydroxy-7, 8, 9, 10-tetrahydro-6, 6,  
110 9-trimethyl-6H-dibenzo -b,d] pyran; Synhexyl);

111 Peyote;

112 N-ethyl-3-piperidyl benzilate;

113 N-methyl-3-piperidyl benzilate;

114 Psilocybin;

115 Psilocyn;

116 Salvinorin A;

117 Tetrahydrocannabinols, except as present in marijuana and dronabinol in sesame oil and encapsulated

- 118 in a soft gelatin capsule in a drug product approved by the U.S. Food and Drug Administration;  
 119 Hashish oil (some trade or other names: hash oil; liquid marijuana; liquid hashish);  
 120 2,5-dimethoxyamphetamine (some trade or other names: 2,5-dimethoxy-a-methylphenethylamine;  
 121 2,5-DMA);  
 122 3,4-methylenedioxymethamphetamine (MDMA), its optical, positional and geometric isomers, salts  
 123 and salts of isomers;  
 124 3,4-methylenedioxymethamphetamine (also known as N-ethyl-alpha-methyl-3,4  
 125 (methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA);  
 126 N-hydroxy-3,4-methylenedioxymethamphetamine (some other names:  
 127 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA);  
 128 4-bromo-2,5-dimethoxyamphetamine (some trade or other names:  
 129 4-bromo-2,5-dimethoxy-a-methylphenethylamine; 4-bromo-2,5-DMA);  
 130 4-methoxyamphetamine (some trade or other names: 4-methoxy-a-methylphenethylamine;  
 131 paramethoxyamphetamine; PMA);  
 132 Ethylamine analog of phencyclidine (some other names: N-ethyl-1-phenylcyclohexylamine,  
 133 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE);  
 134 Pyrrolidine analog of phencyclidine (some other names: 1-(1-phenylcyclohexyl) -pyrrolidine, PCPy,  
 135 PHP);  
 136 Thiophene analog of phencyclidine (some other names: 1-(2-thienyl) -cyclohexyl]-piperidine,  
 137 2-thienyl analog of phencyclidine, TPCP, TCP);  
 138 1-(2-thienyl) Cyclohexyl]pyrrolidine (other name: TCPy);  
 139 3,4-methylenedioxypyrovalerone (other name: MDPV);  
 140 4-methylmethcathinone (other names: mephedrone, 4-MMC);  
 141 3,4-methylenedioxymethcathinone (other name: methylone);  
 142 Naphthylpyrovalerone (other name: naphyrone);  
 143 4-fluoromethcathinone (other name: flephedrone, 4-FMC);  
 144 4-methoxymethcathinone (other names: methedrone; bk-PMMA);  
 145 Ethcathinone (other name: N-ethylcathinone);  
 146 3,4-methylenedioxyethcathinone (other name: ethylone);  
 147 Beta-keto-N-methyl-3,4-benzodioxolylbutanamine (other name: butylone);  
 148 N,N-dimethylcathinone (other name: metamfepramone);  
 149 Alpha-pyrrolidinopropiophenone (other name: alpha-PPP);  
 150 4-methoxy-alpha-pyrrolidinopropiophenone (other name: MOPPP);  
 151 3,4-methylenedioxy-alpha-pyrrolidinopropiophenone (other name: MDPPP);  
 152 Alpha-pyrrolidinovalerophenone (other name: alpha-PVP);  
 153 6,7-dihydro-5H-indeno-(5,6-d)-1,3-dioxol-6-amine (other name: MDAI);  
 154 3-fluoromethcathinone (other name: 3-FMC);  
 155 4-Ethyl-2,5-dimethoxyphenethylamine (other name: 2C-E);  
 156 4-Iodo-2,5-dimethoxyphenethylamine (other name: 2C-I);  
 157 4-Methylethcathinone (other name: 4-MEC);  
 158 4-Ethylmethcathinone (other name: 4-EMC);  
 159 N,N-diallyl-5-methoxytryptamine (other name: 5-MeO-DALT);  
 160 Beta-keto-methylbenzodioxolylpentanamine (other name: Pentylone, bk-MBDP);  
 161 Alpha-methylamino-butyrophenone (other name: Buphedrone);  
 162 Alpha-methylamino-valerophenone (other name: Pentedrone);  
 163 3,4-Dimethylmethcathinone (other name: 3,4-DMMC);  
 164 4-methyl-alpha-pyrrolidinopropiophenone (other name: MPPP);  
 165 4-Iodo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names: 25-I,  
 166 25I-NBOMe);  
 167 Methoxetamine (other names: MXE, 3-MeO-2-Oxo-PCE);  
 168 4-Fluoromethamphetamine (other name: 4-FMA);  
 169 4-Fluoroamphetamine (other name: 4-FA);  
 170 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (other name: 2C-D);  
 171 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (other name: 2C-C);  
 172 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-2);  
 173 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (other name: 2C-T-4);  
 174 2-(2,5-Dimethoxyphenyl)ethanamine (other name: 2C-H);  
 175 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (other name: 2C-N);  
 176 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (other name: 2C-P);  
 177 (2-aminopropyl)benzofuran (other name: APB);  
 178 (2-aminopropyl)-2,3-dihydrobenzofuran (other name: APDB);

- 179 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:  
180 2C-C-NBOMe, 25C-NBOMe);
- 181 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)methyl]-benzeneethanamine (other names:  
182 2C-B-NBOMe, 25B-NBOMe);
- 183 Acetoxymethyltryptamine (other names: AcO-Psilocin, AcO-DMT, Psilacetin);
- 184 Benocyclidine (other names: BCP, BTCP);
- 185 Alpha-pyrrolidinobutylphenone (other name: alpha-PBP);
- 186 3,4-methylenedioxy-N,N-dimethylcathinone (other names: Dimethylone, bk-MDDMA);
- 187 4-bromomethylcathinone (other name: 4-BMC);
- 188 4-chloromethylcathinone (other name: 4-CMC);
- 189 4-Iodo-2,5-dimethoxy-N-[(2-hydroxyphenyl)methyl]-benzeneethanamine (other name: 25I-NBOH);
- 190 Alpha-Pyrrolidinohexylphenone (other name: alpha-PHP);
- 191 Alpha-Pyrrolidinoheptylphenone (other name: PV8).
- 192 4. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture  
193 or preparation which contains any quantity of the following substances having a depressant effect on the  
194 central nervous system, including its salts, isomers and salts of isomers whenever the existence of such  
195 salts, isomers and salts of isomers is possible within the specific chemical designation:
- 196 Gamma hydroxybutyric acid (some other names include GHB; gamma hydroxybutyrate;  
197 4-hydroxybutyrate; 4-hydroxybutanoic acid; sodium oxybate; sodium oxybutyrate);
- 198 Mecloqualone;
- 199 Methaqualone;
- 200 Etizolam.
- 201 5. Unless specifically excepted or unless listed in another schedule, any material, compound, mixture  
202 or preparation which contains any quantity of the following substances having a stimulant effect on the  
203 central nervous system, including its salts, isomers and salts of isomers:
- 204 Aminorex (some trade or other names; aminoxaphen; 2-amino-5-phenyl-2-oxazoline; 4,  
205 5-dihydro-5-phenyl-2-oxazolamine);
- 206 N-Benzylpiperazine (some other names: BZP, 1-benzylpiperazine);
- 207 Fenethylamine;
- 208 Ethylamphetamine;
- 209 Cathinone (some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-aminopropiophenone,  
210 2-aminopropiophenone, norephedrine), and any plant material from which Cathinone may be derived;
- 211 Methcathinone (some other names: 2-(methylamino)-propylphenone;  
212 alpha-(methylamino)-propylphenone; 2-(methylamino)-1-phenylpropan-1-one;  
213 alpha-N-methylaminopropiophenone; monomethylpropion; ephedrine; N-methylcathinone;  
214 methylcathinone; AL-464; AL-422; AL-463 and UR 1432);
- 215 Cis-4-methylaminorex (other name: cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
- 216 N,N-dimethylamphetamine (other names: N,N-alpha-trimethyl-benzeneethanamine,  
217 N,N-alpha-trimethylphenethylamine).
- 218 6. Any material, compound, mixture or preparation containing any quantity of the following  
219 substances:
- 220 N-3-methyl-1-(2-phenethyl)-4-piperidyl-N-phenylpropanamide (other name: 3-methylfentanyl), its  
221 optical and geometric isomers, salts, and salts of isomers;
- 222 1-methyl-4-phenyl-4-propionoxypiperidine (other name: MPPP), its optical isomers, salts and salts of  
223 isomers;
- 224 1-(2-phenylethyl)-4-phenyl-4-acetyloxypiperidine (other name: PEPAP), its optical isomers, salts and  
225 salts of isomers;
- 226 N-1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl propionanilide (other names:  
227 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine), alpha-methylfentanyl);
- 228 N-1-(1-methyl-2-phenethyl)-4-piperidyl-N-phenylacetamide (other name:  
229 acetyl-alpha-methylfentanyl), its optical isomers, salts and salts of isomers;
- 230 N-1-(1-methyl-2-2-thienyl)ethyl-4 piperidyl-N-phenylpropanamide (other name:  
231 alpha-methylthiofentanyl), its optical isomers, salts and salts of isomers;
- 232 N-1-benzyl-4-piperidyl-N-phenylpropanamide (other name: benzylfentanyl), its optical isomers, salts  
233 and salts of isomers;
- 234 N-1-(2-hydroxy-2-phenyl) ethyl-4-piperidyl-N-phenylpropanamide (other name:  
235 beta-hydroxyfentanyl), its optical isomers, salts and salts of isomers;
- 236 N-3-methyl-1-(2-hydroxy-2-phenethyl)-4-piperidyl-N-phenylpropanamide (other name:  
237 beta-hydroxy-3-methylfentanyl), its optical and geometric isomers, salts and salts of isomers;
- 238 N-(3-methyl-1-(2-thienyl)ethyl-4-piperidyl)-N-phenylpropanamide (other name:  
239 3-methylthiofentanyl), its optical and geometric isomers, salts and salts of isomers;

240 N-1-(2-thienyl) Methyl-4-piperidyl]-N-phenylpropanamide (other name: thienylfentanyl), its optical  
 241 isomers, salts and salts of isomers;  
 242 N-phenyl-N-1-(2-thienyl)ethyl-4-piperidiny]-propanamide (other name: thiofentanyl), its optical  
 243 isomers, salts and salts of isomers;  
 244 N-(4-fluorophenyl)-N-1-(2-phenethyl)-4-piperidiny]-propanamide (other name: para-fluorofentanyl),  
 245 its optical isomers, salts and salts of isomers;  
 246 *Acetyl fentanyl (other name: desmethyl fentanyl).*  
 247 7. Any substance that contains one or more cannabimimetic agents or that contains their salts,  
 248 isomers, and salts of isomers whenever the existence of such salts, isomers, and salts of isomers is  
 249 possible within the specific chemical designation, and any preparation, mixture, or substance containing,  
 250 or mixed or infused with, any detectable amount of one or more cannabimimetic agents.  
 251 a. "Cannabimimetic agents" includes any substance that is within any of the following structural  
 252 classes:  
 253 2-(3-hydroxycyclohexyl)phenol with substitution at the 5-position of the phenolic ring by alkyl or  
 254 alkenyl, whether or not substituted on the cyclohexyl ring to any extent;  
 255 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl) Methane with substitution at the nitrogen atom of  
 256 the indole ring, whether or not further substituted on the indole ring to any extent, whether or not  
 257 substituted on the naphthoyl or naphthyl ring to any extent;  
 258 3-(1-naphthoyl)pyrrole with substitution at the nitrogen atom of the pyrrole ring, whether or not  
 259 further substituted in the pyrrole ring to any extent, whether or not substituted on the naphthoyl ring to  
 260 any extent;  
 261 1-(1-naphthylmethyl)indene with substitution of the 3-position of the indene ring, whether or not  
 262 further substituted in the indene ring to any extent, whether or not substituted on the naphthyl ring to  
 263 any extent;  
 264 3-phenylacetylindole or 3-benzoylindole with substitution at the nitrogen atom of the indole ring,  
 265 whether or not further substituted in the indole ring to any extent, whether or not substituted on the  
 266 phenyl ring to any extent;  
 267 3-cyclopropylindole with substitution at the nitrogen atom of the indole ring, whether or not further  
 268 substituted on the indole ring to any extent, whether or not substituted on the cyclopropyl ring to any  
 269 extent;  
 270 3-adamantoylindole with substitution at the nitrogen atom of the indole ring, whether or not further  
 271 substituted on the indole ring to any extent, whether or not substituted on the adamantyl ring to any  
 272 extent;  
 273 N-(adamantyl)-indole-3-carboxamide with substitution at the nitrogen atom of the indole ring,  
 274 whether or not further substituted on the indole ring to any extent, whether or not substituted on the  
 275 adamantyl ring to any extent; and  
 276 N-(adamantyl)-indazole-3-carboxamide with substitution at a nitrogen atom of the indazole ring,  
 277 whether or not further substituted on the indazole ring to any extent, whether or not substituted on the  
 278 adamantyl ring to any extent.  
 279 b. The term "cannabimimetic agents" includes:  
 280 5-(1,1-Dimethylheptyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497);  
 281 5-(1,1-Dimethylhexyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C6 homolog);  
 282 5-(1,1-Dimethyloctyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C8 homolog);  
 283 5-(1,1-Dimethylnonyl)-2-[3-hydroxycyclohexyl]-phenol (other name: CP 47,497 C9 homolog);  
 284 1-pentyl-3-(1-naphthoyl)indole (other names: JWH-018, AM-678);  
 285 1-butyl-3-(1-naphthoyl)indole (other name: JWH-073);  
 286 1-pentyl-3-(2-methoxyphenylacetyl)indole (other name: JWH-250);  
 287 1-hexyl-3-(naphthalen-1-oyl)indole (other name: JWH-019);  
 288 1-[2-(4-morpholinyl)ethyl]-3-(1-naphthoyl)indole (other name: JWH-200);  
 289 (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chr-  
 290 omen-1-ol (other name: HU-210);  
 291 1-pentyl-3-(4-methoxy-1-naphthoyl)indole (other name: JWH-081);  
 292 1-pentyl-3-(4-methyl-1-naphthoyl)indole (other name: JWH-122);  
 293 1-pentyl-3-(2-chlorophenylacetyl)indole (other name: JWH-203);  
 294 1-pentyl-3-(4-ethyl-1-naphthoyl)indole (other name: JWH-210);  
 295 1-pentyl-3-(4-chloro-1-naphthoyl)indole (other name: JWH-398);  
 296 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (other name: AM-694);  
 297 1-((N-methylpiperidin-2-yl) Methyl)-3-(1-naphthoyl)indole (other name: AM-1220);  
 298 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (other name: AM-2201);  
 299 1-[(N-methylpiperidin-2-yl) Methyl]-3-(2-iodobenzoyl)indole (other name: AM-2233);  
 300 Pravadoline (4-methoxyphenyl)-[2-methyl-1-(2-(4-morpholinyl)ethyl)indol-3-yl]methanone (other

301 name: WIN 48,098);  
 302 1-pentyl-3-(4-methoxybenzoyl)indole (other names: RCS-4, SR-19);  
 303 1-(2-cyclohexylethyl)-3-(2-methoxyphenylacetyl)indole (other names: RCS-8, SR-18);  
 304 1-pentyl-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: UR-144);  
 305 1-(5-fluoropentyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: XLR-11);  
 306 N-adamantyl-1-fluoropentylindole-3-carboxamide (other name: STS-135);  
 307 N-adamantyl-1-pentylindazole-3-carboxamide (other name: AKB48);  
 308 1-pentyl-3-(1-adamantoyl)indole (other name: AB-001);  
 309 (8-quinoliny)(1-pentylindol-3-yl)carboxylate (other name: PB-22);  
 310 (8-quinoliny)(1-(5-fluoropentyl) Indol-3-yl)carboxylate (other name: 5-fluoro-PB-22);  
 311 (8-quinoliny)(1-cyclohexylmethyl-indol-3-yl)carboxylate (other name: BB-22);  
 312 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name: AB-PINACA);  
 313 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)indazole-3-carboxamide (other name:  
 314 AB-FUBINACA);  
 315 1-(5-fluoropentyl)-3-(1-naphthoyl)indazole (other name: THJ-2201);  
 316 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentylindazole-3-carboxamide (other name:  
 317 ADB-PINACA);  
 318 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other name:  
 319 AB-CHMINACA);  
 320 N-(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)indazole-3-carboxamide (other name:  
 321 5-fluoro-AB-PINACA);  
 322 *N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)indazole-3-carboxamide (other names:*  
 323 *ADB-CHMINACA, MAB-CHMINACA);*  
 324 *Methyl-2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (other name:*  
 325 *5-fluoro-AMB);*  
 326 *1-naphthalenyl 1-(5-fluoropentyl)-1H-indole-3-carboxylate (other name: NM-2201);*  
 327 *1-(4-fluorobenzyl)-3-(2,2,3,3-tetramethylcyclopropylmethanone)indole (other name: FUB-144);*  
 328 *1-(5-fluoropentyl)-3-(4-methyl-1-naphthoyl)indole (other name MAM-2201).*  
 329 **2. That the provisions of this act may result in a net increase in periods of imprisonment or**  
 330 **commitment. Pursuant to § 30-19.1:4, the estimated amount of the necessary appropriation is \$0**  
 331 **for periods of imprisonment in state adult correctional facilities and cannot be determined for**  
 332 **periods of commitment to the custody of the Department of Juvenile Justice.**