

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101821\  
 Data File : BP007532.D  
 Acq On : 20 Oct 2021 16:59  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 87 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020707

Quant Time: Oct 20 17:36:39 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP101421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 20 10:49:33 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| -----                         |        |      |          |       |        |          |
| Internal Standards            |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.952  | 152  | 186306   | 20.00 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.757 | 136  | 751888   | 20.00 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 473686   | 20.00 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.340 | 188  | 1016636  | 20.00 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.428 | 240  | 1053226  | 20.00 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.874 | 264  | 1120578  | 20.00 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 30856    | 6.54  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 239295   | 18.53 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 275252   | 18.92 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67   | 156078   | 17.99 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 221086   | 19.32 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 215688   | 19.25 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.111  | 128  | 100424   | 21.11 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 103319   | 23.25 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 221717   | 20.45 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.887 | 131  | 285924   | 19.50 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.999 | 166  | 608548   | 18.57 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 784100   | 19.04 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.775 | 143  | 108864   | 20.70 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 528922   | 19.29 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.693 | 200  | 76453    | 18.77 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.439 | 188  | 832757   | 19.29 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.669 | 212  | 972015   | 18.46 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.716 | 264  | 1059779  | 19.01 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |       |        |          |
|                               |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane                | 3.417  | 88   | 36175    | 7.05  | ng/uL  | 85       |
| 5) Pyridine                   | 3.817  | 79   | 237421   | 19.01 | ng/ul  | 98       |
| 6) Benzaldehyde               | 7.087  | 77   | 180529   | 22.08 | ng/ul  | 97       |
| 8) Phenol                     | 7.140  | 94   | 278347   | 18.71 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.369  | 93   | 223177   | 18.59 | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.516  | 128  | 225426   | 19.18 | ng/ul  | 97       |
| 13) 2-Methylphenol            | 8.393  | 108  | 211335   | 18.86 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.493  | 45   | 258610   | 16.90 | ng/ul  | 96       |
| 16) Acetophenone              | 8.775  | 105  | 332052   | 19.37 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.763  | 70   | 161595   | 18.22 | ng/ul  | 93       |
| 18) 4-Methylphenol            | 8.722  | 108  | 229659   | 19.24 | ng/ul  | 97       |
| 19) Hexachloroethane          | 9.040  | 117  | 90045    | 18.30 | ng/ul  | 89       |
| 22) Nitrobenzene              | 9.152  | 77   | 241779   | 19.79 | ng/ul  | 97       |
| 23) Isophorone                | 9.687  | 82   | 455900   | 17.83 | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.863  | 139  | 113644   | 22.06 | ng/ul  | 97       |
| 26) 2,4-Dimethylphenol        | 9.928  | 107  | 253964   | 19.31 | ng/ul  | 96       |
| 27) Bis(2-Chloroethoxy)met... | 10.169 | 93   | 288647   | 18.21 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.405 | 162  | 216197   | 20.29 | ng/ul  | 95       |
| 30) Naphthalene               | 10.810 | 128  | 695971   | 19.03 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.910 | 127  | 295480   | 19.92 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 11.105 | 225  | 150057   | 20.06 | ng/ul  | 98       |
| 34) Caprolactam               | 11.675 | 113  | 67309    | 24.08 | ng/ul  | 92       |
| 35) 4-Chloro-3-methylphenol   | 12.034 | 107  | 225569   | 19.49 | ng/ul  | 97       |
| 36) 2-Methylnaphthalene       | 12.416 | 142  | 469807   | 18.80 | ng/ul  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101821\  
 Data File : BP007532.D  
 Acq On : 20 Oct 2021 16:59  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 87 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020707

Quant Time: Oct 20 17:36:39 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP101421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 20 10:49:33 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| 37) 1-Methylnaphthalene       | 12.634 | 142  | 480836   | 19.01 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.787 | 216  | 273053   | 19.94 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.775 | 237  | 132432   | 15.34 | ng/ul | 98       |
| 41) 2,4,6-Trichlorophenol     | 13.022 | 196  | 170795   | 21.21 | ng/ul | 98       |
| 42) 2,4,5-Trichlorophenol     | 13.093 | 196  | 186364   | 22.23 | ng/ul | 97       |
| 43) 1,1'-Biphenyl             | 13.422 | 154  | 636573   | 18.85 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.463 | 162  | 496480   | 19.09 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.657 | 65   | 128233   | 21.21 | ng/ul | 96       |
| 47) Dimethylphthalate         | 14.046 | 163  | 606170   | 18.62 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.157 | 165  | 114145   | 21.15 | ng/ul | 94       |
| 50) Acenaphthylene            | 14.310 | 152  | 790084   | 19.03 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.481 | 138  | 127579   | 21.21 | ng/ul | 98       |
| 52) Acenaphthene              | 14.651 | 153  | 508730   | 18.92 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.687 | 184  | 47486    | 19.04 | ng/ul | 96       |
| 55) 4-Nitrophenol             | 14.787 | 109  | 95868    | 20.18 | ng/ul | 94       |
| 56) Dibenzofuran              | 14.987 | 168  | 748733   | 19.18 | ng/ul | 99       |
| 57) 2,4-Dinitrotoluene        | 14.945 | 165  | 164758   | 21.73 | ng/ul | 91       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.210 | 232  | 152557   | 21.93 | ng/ul | 97       |
| 59) Diethylphthalate          | 15.410 | 149  | 600969   | 18.52 | ng/ul | 99       |
| 61) Fluorene                  | 15.634 | 166  | 580645   | 19.63 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.634 | 204  | 299199   | 19.86 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.645 | 138  | 130297   | 24.44 | ng/ul | 96       |
| 66) 4,6-Dinitro-2-methylph... | 15.710 | 198  | 78469    | 18.62 | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine    | 15.845 | 169  | 518241   | 18.93 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.528 | 248  | 187579   | 20.48 | ng/ul | 98       |
| 69) Hexachlorobenzene         | 16.645 | 284  | 217825   | 19.77 | ng/ul | 97       |
| 70) Atrazine                  | 16.798 | 200  | 197654   | 19.08 | ng/ul | 96       |
| 71) Pentachlorophenol         | 16.992 | 266  | 130769   | 21.44 | ng/ul | 97       |
| 72) Phenanthrene              | 17.381 | 178  | 970246   | 19.09 | ng/ul | 100      |
| 74) Anthracene                | 17.475 | 178  | 976427   | 19.21 | ng/ul | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.393 | 216  | 285070   | 19.33 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.910 | 250  | 271393   | 19.74 | ng/uL | 99       |
| 77) Carbazole                 | 17.739 | 167  | 899874   | 20.15 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.304 | 149  | 1030221  | 18.78 | ng/ul | 99       |
| 80) Fluoranthene              | 19.345 | 202  | 1206219  | 18.46 | ng/ul | 99       |
| 82) Pyrene                    | 19.698 | 202  | 1221750  | 18.55 | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.575 | 149  | 481484   | 19.91 | ng/ul | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.339 | 252  | 391566   | 19.54 | ng/ul | 98       |
| 85) Benzo(a)anthracene        | 21.410 | 228  | 1218958  | 18.84 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.345 | 149  | 712226   | 19.26 | ng/ul | 100      |
| 87) Chrysene                  | 21.463 | 228  | 1172959  | 18.71 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.286 | 149  | 1216894  | 19.25 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.127 | 252  | 1265146  | 18.55 | ng/ul | 98       |
| 91) Benzo(k)fluoranthene      | 23.174 | 252  | 1192518  | 18.75 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 23.769 | 252  | 1224667  | 18.77 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.433 | 276  | 1426649  | 18.88 | ng/ul | 100      |
| 95) Dibenzo(a,h)anthracene    | 26.457 | 278  | 1210548  | 18.81 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 27.215 | 276  | 1217185  | 18.70 | ng/ul | 97       |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quant Time: Oct 20 17:36:39 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP101421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 20 10:49:33 2021  
Response via : Initial Calibration

