

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101823\  
 Data File : BP017715.D  
 Acq On : 18 Oct 2023 18:24  
 Operator : MA/JU  
 Sample : SST00534  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST005645

Quant Time: Oct 19 01:12:37 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101823.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 19 00:53:23 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.887  | 152  | 130752   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.716 | 136  | 480662   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 311914   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.369 | 188  | 692177   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.510 | 240  | 658928   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.057 | 264  | 742733   | 20.000 | ng/u1  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 6843     | 1.952  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 0.000  | 99   | 0d       | 0.000  | ng/u1  |          |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000  | 67   | 0d       | 0.000  | ng/u1  |          |
| 11) 2-Chlorophenol-d4         | 7.410  | 132  | 38946    | 4.191  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 0.000  | 113  | 0d       | 0.000  | ng/u1  |          |
| 21) Nitrobenzene-d5           | 9.093  | 128  | 16271    | 4.060  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 15831    | 3.523  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.346 | 165  | 32999    | 3.876  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 0.000  | 131  | 0d       | 0.000  | ng/u1  |          |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 122369   | 4.575  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.275 | 160  | 127740   | 4.402  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 15.587 | 176  | 104720   | 4.585  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/u1  |          |
| 73) Anthracene-d10            | 17.469 | 188  | 160131   | 4.498  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.727 | 212  | 188479   | 4.713  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.892 | 264  | 181107   | 4.334  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.299  | 88   | 7071     | 1.875  | ng/uL  | 96       |
| 12) 2-Chlorophenol            | 7.446  | 128  | 38828    | 4.192  | ng/u1  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.710  | 70   | 27382    | 3.714  | ng/u1  | 95       |
| 19) Hexachloroethane          | 8.946  | 117  | 19695    | 4.738  | ng/u1  | 93       |
| 22) Nitrobenzene              | 9.140  | 77   | 42400    | 4.148  | ng/u1  | 99       |
| 23) Isophorone                | 9.652  | 82   | 78578    | 4.287  | ng/u1  | 98       |
| 25) 2-Nitrophenol             | 9.846  | 139  | 17683    | 3.696  | ng/u1  | 97       |
| 26) 2,4-Dimethylphenol        | 9.887  | 107  | 39904    | 4.223  | ng/u1  | 97       |
| 27) Bis(2-Chloroethoxy)met... | 10.140 | 93   | 47964    | 4.240  | ng/u1  | 97       |
| 29) 2,4-Dichlorophenol        | 10.375 | 162  | 34382    | 4.207  | ng/u1  | 94       |
| 30) Naphthalene               | 10.769 | 128  | 130813   | 4.697  | ng/u1  | 98       |
| 33) Hexachlorobutadiene       | 10.993 | 225  | 31844    | 4.983  | ng/u1  | 93       |
| 35) 4-Chloro-3-methylphenol   | 12.034 | 107  | 30480    | 3.610  | ng/u1  | 98       |
| 36) 2-Methylnaphthalene       | 12.387 | 142  | 86162    | 4.531  | ng/u1  | 97       |
| 37) 1-Methylnaphthalene       | 12.610 | 142  | 88565    | 4.570  | ng/u1  | 93       |
| 39) 1,2,4,5-Tetrachloroben... | 12.740 | 216  | 53445    | 4.677  | ng/u1  | 98       |
| 41) 2,4,6-Trichlorophenol     | 13.004 | 196  | 27828    | 3.991  | ng/u1  | 96       |
| 42) 2,4,5-Trichlorophenol     | 13.087 | 196  | 29014    | 3.953  | ng/u1  | 94       |
| 43) 1,1'-Biphenyl             | 13.404 | 154  | 116721   | 4.554  | ng/u1  | 99       |
| 44) 2-Chloronaphthalene       | 13.457 | 162  | 95983    | 4.677  | ng/u1  | 98       |
| 45) 2-Nitroaniline            | 13.710 | 65   | 16478    | 3.020  | ng/u1  | 96       |
| 47) Dimethylphthalate         | 14.045 | 163  | 122142   | 4.621  | ng/u1  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.198 | 165  | 18211    | 3.757  | ng/u1  | 88       |
| 50) Acenaphthylene            | 14.310 | 152  | 145905   | 4.442  | ng/u1  | 99       |

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 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD005645

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 19 00:53:23 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| 52) Acenaphthene              | 14.645 | 153  | 106221   | 4.846 | ng/ul | 100      |
| 56) Dibenzofuran              | 14.987 | 168  | 144781   | 4.616 | ng/ul | 96       |
| 57) 2,4-Dinitrotoluene        | 14.992 | 165  | 26407    | 3.532 | ng/ul | 92       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.210 | 232  | 27259    | 4.068 | ng/ul | 94       |
| 59) Diethylphthalate          | 15.404 | 149  | 116060   | 4.414 | ng/ul | 97       |
| 61) Fluorene                  | 15.639 | 166  | 121051   | 4.699 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.634 | 204  | 64379    | 4.704 | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine    | 15.863 | 169  | 94448    | 4.566 | ng/ul | 98       |
| 68) 4-Bromophenyl-phenylether | 16.545 | 248  | 36978    | 4.461 | ng/ul | 97       |
| 69) Hexachlorobenzene         | 16.622 | 284  | 43107    | 4.396 | ng/ul | 96       |
| 72) Phenanthrene              | 17.416 | 178  | 191757   | 4.712 | ng/ul | 99       |
| 74) Anthracene                | 17.510 | 178  | 190424   | 4.633 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.363 | 216  | 52558    | 4.749 | ng/uL | 97       |
| 76) Pentachlorobenzene        | 14.875 | 250  | 52757    | 4.704 | ng/uL | 97       |
| 78) Di-n-butylphthalate       | 18.310 | 149  | 170610   | 3.809 | ng/ul | 99       |
| 80) Fluoranthene              | 19.398 | 202  | 214099   | 4.674 | ng/ul | 100      |
| 82) Pyrene                    | 19.757 | 202  | 234070   | 4.879 | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.627 | 149  | 69440    | 3.819 | ng/ul | 97       |
| 85) Benzo(a)anthracene        | 21.492 | 228  | 228892   | 4.526 | ng/ul | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.374 | 149  | 94159    | 3.457 | ng/ul | 99       |
| 87) Chrysene                  | 21.551 | 228  | 222225   | 4.737 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.268 | 252  | 225946   | 4.443 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.316 | 252  | 231019   | 4.584 | ng/ul | 97       |
| 93) Benzo(a)pyrene            | 23.945 | 252  | 206070   | 4.368 | ng/ul | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.768 | 276  | 251039   | 4.354 | ng/ul | 98       |
| 95) Dibenzo(a,h)anthracene    | 26.798 | 278  | 208640   | 4.331 | ng/ul | 97       |
| 96) Benzo(g,h,i)perylene      | 27.609 | 276  | 207403   | 4.523 | ng/ul | 99       |

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

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