

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120723\  
 Data File : BP018700.D  
 Acq On : 08 Dec 2023 07:01  
 Operator : MA/JU  
 Sample : 05497-02  
 Misc :  
 ALS Vial : 36 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AB1

Quant Time: Dec 08 07:29:20 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 04 21:39:40 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.846  | 152  | 369378   | 20.000 | ng/u1  | -0.01    |
| 20) Naphthalene-d8            | 10.640 | 136  | 1008572  | 20.000 | ng/u1  | -0.01    |
| 38) Acenaphthene-d10          | 14.475 | 164  | 634570   | 20.000 | ng/u1  | -0.01    |
| 64) Phenanthrene-d10          | 17.222 | 188  | 1517248  | 20.000 | ng/u1  | # 0.00   |
| 79) Chrysene-d12              | 21.310 | 240  | 1730674  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.680 | 264  | 2059689  | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 13729    | 1.267  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 50679    | 1.745  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.016  | 99   | 132154   | 3.565  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 149708   | 6.814  | ng/u1  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.375  | 132  | 140923   | 4.872  | ng/u1  | -0.01    |
| 15) 4-Methylphenol-d8         | 8.557  | 113  | 19863    | 0.662  | ng/u1  | -0.01    |
| 21) Nitrobenzene-d5           | 9.004  | 128  | 89889    | 10.014 | ng/u1  | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.728  | 143  | 62590    | 6.725  | ng/u1  | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.263 | 165  | 94138    | 4.947  | ng/u1  | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.781 | 131  | 70994    | 2.855  | ng/u1  | -0.01    |
| 46) Dimethylphthalate-d6      | 13.892 | 166  | 410112   | 7.241  | ng/u1  | -0.01    |
| 49) Acenaphthylene-d8         | 14.169 | 160  | 449192   | 7.203  | ng/u1  | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.681 | 143  | 6283     | 0.676  | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.469 | 176  | 368097   | 7.743  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.586 | 200  | 7641     | 0.840  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.322 | 188  | 570170   | 7.148  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.557 | 212  | 823137   | 6.110  | ng/u1  | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.521 | 264  | 806250   | 6.713  | ng/u1  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 8) Phenol                     | 7.046  | 94   | 41744    | 1.148  | ng/u1  | 96       |
| 16) Acetophenone              | 8.669  | 105  | 213719   | 4.780  | ng/u1  | 99       |
| 72) Phenanthrene              | 17.263 | 178  | 114202   | 1.246  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.233 | 202  | 245005   | 1.539  | ng/u1  | 97       |
| 82) Pyrene                    | 19.586 | 202  | 211078   | 1.314  | ng/u1# | 92       |
| 90) Benzo(b)fluoranthene      | 22.957 | 252  | 166548   | 1.097  | ng/u1# | 94       |
| -----                         |        |      |          |        |        |          |

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

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