

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120123\  
 Data File : BP018494.D  
 Acq On : 02 Dec 2023 01:18  
 Operator : MA/JU  
 Sample : 05332-06  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 YCG99

Quant Time: Dec 02 01:46:45 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 01 21:56:43 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.863  | 152  | 294485   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.663 | 136  | 1238341  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.492 | 164  | 793245   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.239 | 188  | 1595763  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.321 | 240  | 1143574  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.692 | 264  | 1287272  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 33443    | 3.870  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 194601   | 8.407  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 158949   | 5.379  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.199  | 67   | 529627   | 30.238 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.393  | 132  | 480905   | 20.855 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 310420   | 12.979 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.022  | 128  | 329225   | 29.871 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.746  | 143  | 318253   | 27.851 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.281 | 165  | 573649   | 24.551 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.798 | 131  | 683376   | 22.379 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.904 | 166  | 2239679  | 31.635 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 2424943  | 31.106 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.686 | 143  | 44289    | 3.811  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.481 | 176  | 1887534  | 31.762 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200  | 229957   | 24.033 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.339 | 188  | 2845921  | 33.924 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.569 | 212  | 2885941  | 32.420 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.545 | 264  | 2544888  | 33.906 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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