

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090123\  
 Data File : BP016916.D  
 Acq On : 01 Sep 2023 19:54  
 Operator : MA/JU  
 Sample : 04123-02  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH923

Quant Time: Sep 01 23:41:48 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 30 02:25:00 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.016  | 152  | 367919   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.846 | 136  | 1616334  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.675 | 164  | 1036938  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.445 | 188  | 2269172  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.545 | 240  | 1909590  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.110 | 264  | 1840856  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.375  | 96   | 50375    | 5.659  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.811  | 84   | 348907   | 12.826 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.169  | 99   | 262102   | 8.012  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.357  | 67   | 690987   | 33.139 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.540  | 132  | 663122   | 26.959 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.728  | 113  | 469267   | 17.571 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.216  | 128  | 431172   | 35.281 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.934  | 143  | 470338   | 36.512 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.457 | 165  | 770733   | 32.093 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.004 | 131  | 1077696  | 29.262 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.087 | 166  | 2678443  | 34.251 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 3049375  | 34.495 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.886 | 143  | 105448   | 6.524  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.669 | 176  | 2308891  | 35.061 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 405843   | 30.032 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.545 | 188  | 3693436  | 36.721 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.780 | 212  | 4192393  | 39.974 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.945 | 264  | 3363026  | 36.471 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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