

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090123\  
 Data File : BP016938.D  
 Acq On : 02 Sep 2023 08:50  
 Operator : MA/JU  
 Sample : 04209-20  
 Misc :  
 ALS Vial : 36 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH488

Quant Time: Sep 04 00:42:52 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  | 236250   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.846 | 136  | 1108080  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.669 | 164  | 749067   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.439 | 188  | 1704127  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 1543444  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.104 | 264  | 1570001  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 30120    | 5.269  | ng/uL | -0.01    |
| 4) Pyridine-d5                | 3.805  | 84   | 261089   | 14.946 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.164  | 99   | 193713   | 9.221  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.352  | 67   | 612911   | 45.777 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 537592   | 34.036 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.722  | 113  | 378276   | 22.058 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.210  | 128  | 376890   | 44.985 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.928  | 143  | 403148   | 45.650 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.452 | 165  | 668071   | 40.578 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.999 | 131  | 985184   | 39.020 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 14.081 | 166  | 2589862  | 45.846 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.369 | 160  | 2845494  | 44.558 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.881 | 143  | 80240    | 6.872  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.663 | 176  | 2205443  | 46.361 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.787 | 200  | 380314   | 37.474 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.539 | 188  | 3660351  | 48.458 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.775 | 212  | 4302955  | 50.761 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.933 | 264  | 3825112  | 48.639 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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