

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100423\  
 Data File : BP017577.D  
 Acq On : 05 Oct 2023 21:39  
 Operator : MA/JU  
 Sample : 04607-02  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BBJ35

Quant Time: Oct 06 02:03:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 04 16:33:47 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.934  | 152  | 88558    | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8        | 10.775 | 136  | 366933   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.628 | 164  | 235835   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10      | 17.416 | 188  | 559934   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12          | 21.551 | 240  | 594794   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 24.133 | 264  | 676933   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.287  | 96  | 18596   | 8.395  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.734  | 84  | 31601   | 5.450  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.093  | 99  | 60563   | 8.128  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67  | 193601  | 42.582 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.458  | 132 | 190875  | 32.534 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.657  | 113 | 121053  | 20.294 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.146  | 128 | 123328  | 43.688 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.863  | 143 | 131696  | 41.363 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.393 | 165 | 227702  | 38.576 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.946 | 131 | 301462  | 35.852 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 14.045 | 166 | 823812  | 44.783 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.328 | 160 | 904867  | 44.608 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.887 | 143 | 21177   | 6.574  | ng/u1 | 0.01  |
| 60) Fluorene-d10              | 15.634 | 176 | 760359  | 48.730 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.775 | 200 | 122179  | 35.000 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.516 | 188 | 1376820 | 53.121 | ng/u1 | -0.01 |
| 81) Pyrene-d10                | 19.775 | 212 | 1758052 | 53.192 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.963 | 264 | 1833623 | 53.674 | ng/u1 | -0.01 |

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

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

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