

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051023\  
 Data File : BP015055.D  
 Acq On : 10 May 2023 22:12  
 Operator : CG/JU  
 Sample : 02591-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JREB4

Quant Time: May 10 23:26:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP050323.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 10 15:17:56 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.152  | 152  | 298354   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.987 | 136  | 1263385  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.793 | 164  | 764590   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.551 | 188  | 1641339  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.628 | 240  | 1043421  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.216 | 264  | 1222871  | 20.000 | ng/u1 | 0.01     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.452  | 96   | 17020    | 2.121  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1 |          |
| 7) Phenol-d5                  | 7.305  | 99   | 377941   | 14.611 | ng/u1 | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.469  | 67   | 259466   | 16.229 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.675  | 132  | 330063   | 17.354 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.863  | 113  | 320714   | 16.089 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.328  | 128  | 182051   | 20.261 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.058 | 143  | 169970   | 18.422 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.605 | 165  | 338229   | 18.670 | ng/u1 | 0.01     |
| 31) 4-Chloroaniline-d4        | 11.122 | 131  | 66131    | 2.524  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.193 | 166  | 884717   | 16.349 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.487 | 160  | 1263584  | 19.470 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.993 | 143  | 48944    | 5.199  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.787 | 176  | 1068821  | 22.958 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.904 | 200  | 23227    | 2.514  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.645 | 188  | 1477489  | 20.034 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.869 | 212  | 1371234  | 19.485 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.045 | 264  | 1128514  | 18.940 | ng/u1 | 0.01     |

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

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