

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051623\  
 Data File : BP015111.D  
 Acq On : 16 May 2023 16:19  
 Operator : CG/JU  
 Sample : 02768-12  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CONP0

Quant Time: May 17 01:05:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 16 01:39:21 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.134  | 152  | 317626   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.963 | 136  | 1454199  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.775 | 164  | 944381   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.528 | 188  | 2066431  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.604 | 240  | 1709283  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.174 | 264  | 1648460  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 41811    | 5.115  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.881  | 84   | 179187   | 7.733  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.281  | 99   | 228622   | 7.752  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.452  | 67   | 618805   | 36.331 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.658  | 132  | 602983   | 27.809 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.840  | 113  | 437740   | 18.628 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.311  | 128  | 397374   | 34.775 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.034 | 143  | 424587   | 33.394 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.575 | 165  | 703937   | 30.682 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.099 | 131  | 944472   | 27.823 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.175 | 166  | 2763702  | 38.728 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.469 | 160  | 3072662  | 37.647 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.957 | 143  | 93360    | 6.737  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.763 | 176  | 2391447  | 39.714 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.881 | 200  | 448710   | 34.456 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.628 | 188  | 4015201  | 42.434 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.845 | 212  | 4515368  | 45.548 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.010 | 264  | 3678243  | 44.949 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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