

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051623\  
 Data File : BP015101.D  
 Acq On : 16 May 2023 09:55  
 Operator : CG/JU  
 Sample : PB152799BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK799

Quant Time: May 17 00:34:06 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  | 302883   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.963 | 136  | 1414911  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.775 | 164  | 926458   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.534 | 188  | 2049460  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.610 | 240  | 1744389  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.186 | 264  | 1780135  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 51132    | 6.559  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.870  | 84   | 733380   | 33.190 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.281  | 99   | 890854   | 31.678 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.452  | 67   | 573200   | 35.292 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.658  | 132  | 711548   | 34.413 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.846  | 113  | 763714   | 34.081 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.310  | 128  | 368403   | 33.135 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.040 | 143  | 410032   | 33.144 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.581 | 165  | 713257   | 31.951 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.099 | 131  | 1132009  | 34.273 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.181 | 166  | 2468867  | 35.265 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.475 | 160  | 2787518  | 34.814 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.963 | 143  | 413823   | 30.441 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.769 | 176  | 2071071  | 35.059 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.887 | 200  | 348972   | 27.019 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.634 | 188  | 3283040  | 34.983 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.851 | 212  | 3748572  | 37.052 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.016 | 264  | 3299345  | 37.337 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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