

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110822\  
 Data File : BP012567.D  
 Acq On : 09 Nov 2022 00:23  
 Operator : CG/JU  
 Sample : N5407-17  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW4Z4

Quant Time: Nov 09 00:50:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 08 22:40:23 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.769  | 152  | 217045   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.569 | 136  | 1076136  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.422 | 164  | 768097   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.186 | 188  | 1737170  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.280 | 240  | 1616098  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.656 | 264  | 1473497  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.228  | 96   | 19686    | 3.652 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.663  | 84   | 64668    | 4.176 ng/u1  | 0.01     |
| 7) Phenol-d5                  | 6.963  | 99   | 146757   | 7.539 ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.110  | 67   | 352574   | 30.339 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.304  | 132  | 361595   | 25.435 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.498  | 113  | 283271   | 18.398 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 251078   | 35.979 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 258399   | 35.675 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.204 | 165  | 462609   | 29.471 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.722 | 131  | 342163   | 14.639 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.839 | 166  | 1878274  | 35.722 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.116 | 160  | 1968332  | 32.757 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.698 | 143  | 45642    | 4.586 ng/u1  | 0.02     |
| 60) Fluorene-d10              | 15.422 | 176  | 1627085  | 36.143 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 287123   | 31.746 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.286 | 188  | 2824386  | 37.734 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.527 | 212  | 3328792  | 41.026 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.503 | 264  | 2828740  | 39.290 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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