

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092424\  
 Data File : BP022042.D  
 Acq On : 26 Sep 2024 00:55  
 Operator : RC/JU  
 Sample : PB163459BL  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK459

Quant Time: Sep 26 03:19:49 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 24 15:23:13 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 101646   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.463 | 136  | 381135   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.310 | 164  | 282984   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.110 | 188  | 694653   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 796113   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.827 | 264  | 875161   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.223  | 96   | 18211    | 7.018  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.634  | 84   | 231610   | 31.224 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 271899   | 32.701 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.046  | 67   | 170804   | 33.361 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 206301   | 34.293 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.405  | 113  | 216069   | 32.794 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.846  | 128  | 101879   | 35.714 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 118660   | 36.686 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 240302   | 34.211 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.599 | 131  | 289173   | 34.503 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.722 | 166  | 808600   | 36.965 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.998 | 160  | 845727   | 36.516 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.516 | 143  | 113511   | 30.337 | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.322 | 176  | 700011   | 36.177 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.451 | 200  | 125228   | 29.037 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.210 | 188  | 1229709  | 37.920 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 1576361  | 38.219 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.604 | 264  | 1701144  | 36.957 | ng/u1 | 0.00     |

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

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

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