

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071024\  
 Data File : BP020891.D  
 Acq On : 10 Jul 2024 22:31  
 Operator : MA/JU  
 Sample : PB161824BL  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Jul 11 00:13:53 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 10 18:38:19 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.722  | 152  | 208948   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.481 | 136  | 865337   | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10      | 14.351 | 164  | 571019   | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10      | 17.174 | 188  | 1184849  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.621 | 240  | 1055491  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 24.968 | 264  | 1225470  | 20.000 | ng/ul | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.257  | 96  | 27022   | 5.885  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.663  | 84  | 347788  | 26.163 | ng/ul | 0.00  |
| 7) Phenol-d5                  | 6.916  | 99  | 466576  | 27.389 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.069  | 67  | 297365  | 28.804 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.263  | 132 | 396220  | 28.230 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8         | 8.428  | 113 | 383882  | 27.133 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5           | 8.869  | 128 | 198479  | 29.530 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.586  | 143 | 230468  | 29.959 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.116 | 165 | 374287  | 26.907 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.633 | 131 | 433154  | 23.505 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6      | 13.757 | 166 | 1271265 | 30.625 | ng/ul | -0.01 |
| 49) Acenaphthylene-d8         | 14.039 | 160 | 1375221 | 29.505 | ng/ul | -0.01 |
| 54) 4-Nitrophenol-d4          | 14.610 | 143 | 156775  | 26.545 | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.374 | 176 | 1057237 | 31.045 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.516 | 200 | 185690  | 25.902 | ng/ul | 0.00  |
| 73) Anthracene-d10            | 17.274 | 188 | 1590626 | 30.226 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.657 | 212 | 1875706 | 28.744 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.739 | 264 | 1837727 | 30.406 | ng/ul | -0.02 |

Target Compounds Qvalue

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

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