

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051424\  
 Data File : BP020370.D  
 Acq On : 14 May 2024 16:54  
 Operator : MA/JU  
 Sample : P2448-10  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SVOC-GPC2-BLANK

Quant Time: May 14 23:59:40 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP050124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 14 23:58:44 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.593  | 152  | 64128    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.363 | 136  | 238520   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.275 | 164  | 181858   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.116 | 188  | 440607   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.610 | 240  | 467466   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.957 | 264  | 531715   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 12       | 0.008  | ng/uL | 0.09     |
| 4) Pyridine-d5                | 3.593  | 84   | 192      | 0.044  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.787  | 99   | 2        | 0.000  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.952  | 67   | 49       | 0.015  | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.134  | 132  | 7        | 0.002  | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.316  | 113  | 2        | 0.000  | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.757  | 128  | 10       | 0.006  | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.481  | 143  | 5        | 0.002  | ng/ul | 0.01     |
| 28) 2,4-Dichlorophenol-d3     | 10.075 | 165  | 33       | 0.009  | ng/ul | 0.07     |
| 31) 4-Chloroaniline-d4        | 10.522 | 131  | 15       | 0.003  | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.687 | 166  | 17       | 0.001  | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 13.975 | 160  | 852      | 0.058  | ng/ul | 0.02     |
| 54) 4-Nitrophenol-d4          | 14.545 | 143  | 7        | 0.003  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.304 | 176  | 3        | 0.000  | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 4        | 0.001  | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.286 | 188  | 4050     | 0.212  | ng/ul | 0.07     |
| 81) Pyrene-d10                | 19.604 | 212  | 5        | 0.000  | ng/ul | -0.01    |
| 92) Benzo(a)pyrene-d12        | 24.721 | 264  | 599      | 0.023  | ng/ul | 0.00     |

Target Compounds Qvalue

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

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