

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040523\  
 Data File : BP014559.D  
 Acq On : 05 Apr 2023 15:57  
 Operator : CG/JU  
 Sample : 02122-13  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CB9C5

Quant Time: Apr 06 00:57:48 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 06 00:55:13 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.987  | 152  | 117652   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.810 | 136  | 512263   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.640 | 164  | 331624   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.398 | 188  | 715866   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 627433   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 693910   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 14266    | 4.702  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.811  | 84   | 164      | 0.021  | ng/ul | 0.01     |
| 7) Phenol-d5                  | 7.169  | 99   | 52369    | 4.826  | ng/ul | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.316  | 67   | 193431   | 27.785 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.522  | 132  | 164791   | 21.013 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 107280   | 12.257 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.169  | 128  | 106115   | 25.660 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.893  | 143  | 106589   | 22.509 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.440 | 165  | 176803   | 20.501 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.963 | 131  | 172057   | 15.146 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 835967   | 31.090 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.334 | 160  | 868347   | 29.104 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.851 | 143  | 9        | 0.002  | ng/ul | -0.01    |
| 60) Fluorene-d10              | 15.634 | 176  | 697535   | 31.108 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.769 | 200  | 86565    | 16.434 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.504 | 188  | 1109322  | 32.020 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.733 | 212  | 1326039  | 31.464 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.827 | 264  | 1225635  | 33.245 | ng/ul | 0.00     |

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

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