

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082823\  
 Data File : BP016854.D  
 Acq On : 30 Aug 2023 04:08  
 Operator : MA/JU  
 Sample : 04148-11  
 Misc :  
 ALS Vial : 73 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BBK07

Quant Time: Aug 30 04:48:52 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 30 02:25:00 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 418871   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.852 | 136  | 1856415  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.675 | 164  | 1160094  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.445 | 188  | 2494697  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 1864704  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.110 | 264  | 1695807  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 49986    | 4.932  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.817  | 84   | 241141   | 7.786  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.163  | 99   | 231845   | 6.225  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.363  | 67   | 814351   | 34.305 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.540  | 132  | 718624   | 25.661 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.728  | 113  | 465276   | 15.302 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.222  | 128  | 502222   | 35.780 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.940  | 143  | 510573   | 34.509 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.463 | 165  | 873357   | 31.663 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.010 | 131  | 536521   | 12.684 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.092 | 166  | 3388654  | 38.733 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 3727877  | 37.693 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.887 | 143  | 67777    | 3.748  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.675 | 176  | 2966903  | 40.271 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 426061   | 28.678 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.545 | 188  | 4840180  | 43.771 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.780 | 212  | 5415133  | 52.876 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.939 | 264  | 3814326  | 44.904 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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