

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082823\  
 Data File : BP016791.D  
 Acq On : 28 Aug 2023 14:14  
 Operator : MA/JU  
 Sample : 04134-12  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCHK6

Quant Time: Aug 29 00:48:40 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 25 01:47:30 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.028  | 152  | 357583   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8            | 10.857 | 136  | 1568015  | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10          | 14.687 | 164  | 987566   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.451 | 188  | 2123174  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.551 | 240  | 1384587  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.121 | 264  | 1297237  | 20.000 | ng/ul | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 49785    | 5.754  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 316598   | 11.974 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.169  | 99   | 243875   | 7.670  | ng/ul | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67   | 678843   | 33.498 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 645895   | 27.017 | ng/ul | -0.01    |
| 15) 4-Methylphenol-d8         | 8.734  | 113  | 439845   | 16.945 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.228  | 128  | 421167   | 35.524 | ng/ul | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.946  | 143  | 430753   | 34.469 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.469 | 165  | 717488   | 30.797 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.016 | 131  | 1027290  | 28.753 | ng/ul | -0.01    |
| 46) Dimethylphthalate-d6      | 14.098 | 166  | 2700516  | 36.260 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.387 | 160  | 3048216  | 36.205 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143  | 75615    | 4.912  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.681 | 176  | 2460817  | 39.237 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.804 | 200  | 361090   | 28.558 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.557 | 188  | 4197440  | 44.601 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.786 | 212  | 4373736  | 57.516 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.951 | 264  | 2990621  | 46.024 | ng/ul | -0.01    |

Target Compounds Qvalue

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

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