

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041923\  
 Data File : BP014727.D  
 Acq On : 19 Apr 2023 13:21  
 Operator : CG/JU  
 Sample : 02378-04  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 COMJ7

Quant Time: Apr 20 01:46:13 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP041823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 19 01:37:13 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.152  | 152  | 178372   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.981 | 136  | 743653   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.787 | 164  | 471450   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.534 | 188  | 1073875  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.610 | 240  | 1051927  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.186 | 264  | 1191745  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.487  | 96   | 20752    | 4.624  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.922  | 84   | 51976    | 4.254  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.287  | 99   | 82793    | 5.368  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.469  | 67   | 339674   | 35.701 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.675  | 132  | 289199   | 25.250 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.846  | 113  | 127350   | 10.592 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.322  | 128  | 181656   | 38.983 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.052 | 143  | 158268   | 37.604 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.587 | 165  | 338054   | 29.582 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.110 | 131  | 207084   | 11.873 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.187 | 166  | 1546496  | 41.428 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.481 | 160  | 1681124  | 38.860 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.945 | 143  | 27271    | 4.593  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.775 | 176  | 1333730  | 41.638 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.881 | 200  | 114552   | 26.359 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.634 | 188  | 2044997  | 39.419 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.845 | 212  | 2552900  | 42.365 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.021 | 264  | 2719166  | 44.023 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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