

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032923\  
 Data File : BP014366.D  
 Acq On : 29 Mar 2023 16:30  
 Operator : CG/JU  
 Sample : 01982-22  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EYGC9

Quant Time: Mar 30 01:03:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 30 00:56:20 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 139591   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.846 | 136  | 616685   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.675 | 164  | 425867   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.433 | 188  | 976371   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.521 | 240  | 918113   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.045 | 264  | 864379   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 15061    | 4.184  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.828  | 84   | 179289   | 19.436 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.181  | 99   | 252477   | 19.612 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.346  | 67   | 175236   | 21.215 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 193758   | 20.824 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.734  | 113  | 204350   | 19.679 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.199  | 128  | 98498    | 19.785 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.922  | 143  | 115957   | 20.341 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.469 | 165  | 190360   | 18.336 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.993 | 131  | 265905   | 19.444 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.081 | 166  | 740491   | 21.445 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.369 | 160  | 799615   | 20.870 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.898 | 143  | 82524    | 15.820 | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.669 | 176  | 646677   | 22.458 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 96730    | 13.464 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.533 | 188  | 1032510  | 21.851 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 1195484  | 19.385 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.880 | 264  | 1023763  | 22.293 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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