

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051523\  
 Data File : BP015084.D  
 Acq On : 15 May 2023 20:49  
 Operator : CG/JU  
 Sample : 02592-19  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JREE0

Quant Time: May 16 01:41:59 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 16 01:39:21 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.134  | 152  | 308571   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.969 | 136  | 1406088  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.775 | 164  | 908735   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.528 | 188  | 2009932  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.610 | 240  | 1747209  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.180 | 264  | 1843032  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 50995    | 6.421  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.875  | 84   | 780738   | 34.681 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.287  | 99   | 1560237  | 54.458 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.457  | 67   | 994239   | 60.086 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.663  | 132  | 1255725  | 59.611 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.846  | 113  | 1410107  | 61.767 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.316  | 128  | 672810   | 60.894 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.040 | 143  | 748177   | 60.857 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.581 | 165  | 1315176  | 59.284 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.104 | 131  | 1995267  | 60.788 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.181 | 166  | 4488456  | 65.364 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.475 | 160  | 4958311  | 63.133 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.963 | 143  | 338995   | 25.423 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.769 | 176  | 3735792  | 64.473 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.886 | 200  | 306946   | 24.232 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.633 | 188  | 5893781  | 64.038 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.851 | 212  | 6633394  | 65.460 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.021 | 264  | 6406095  | 70.020 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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