

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP050923\  
 Data File : BP015027.D  
 Acq On : 09 May 2023 20:20  
 Operator : CG/JU  
 Sample : 02694-08  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JREM8

Quant Time: May 09 22:15:38 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP050323.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri May 05 12:12:36 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.157  | 152  | 354014   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.993 | 136  | 1678001  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.804 | 164  | 1087921  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.557 | 188  | 2391721  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.639 | 240  | 1926656  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.233 | 264  | 1665170  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.452  | 96   | 34261    | 3.598  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.899  | 84   | 148245   | 5.696  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.299  | 99   | 176597   | 5.754  | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.475  | 67   | 499660   | 26.339 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.681  | 132  | 511918   | 22.683 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.863  | 113  | 348364   | 14.729 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.334  | 128  | 340878   | 28.563 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.063 | 143  | 369628   | 30.162 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.604 | 165  | 606335   | 25.199 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.128 | 131  | 768161   | 22.071 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.204 | 166  | 2270209  | 29.484 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.498 | 160  | 2543686  | 27.546 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.987 | 143  | 63055    | 4.707  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.792 | 176  | 1930502  | 29.142 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.910 | 200  | 316826   | 23.532 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.657 | 188  | 3215117  | 29.918 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.875 | 212  | 3620115  | 27.859 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.063 | 264  | 2636293  | 32.493 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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