

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102521\  
 Data File : BP007715.D  
 Acq On : 27 Oct 2021 12:35  
 Operator : CG/JU  
 Sample : M4304-08  
 Misc :  
 ALS Vial : 72 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 P067-TW001-01

Quant Time: Oct 27 13:39:23 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP102521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 25 16:57:16 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 183261   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.740 | 136  | 738649   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.569 | 164  | 470708   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.328 | 188  | 1014060  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.416 | 240  | 1092202  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.862 | 264  | 1147115  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.369  | 96   | 25515    | 6.325  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.787  | 84   | 138675   | 11.367 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 107093   | 7.694  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.263  | 67   | 303010   | 39.208 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 324454   | 29.970 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.640  | 113  | 198205   | 18.185 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.093  | 128  | 193243   | 38.390 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 193140   | 35.131 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.357 | 165  | 359169   | 32.096 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.869 | 131  | 488673   | 33.026 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.981 | 166  | 1262355  | 39.592 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.263 | 160  | 1462732  | 37.430 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 36981    | 6.327  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.569 | 176  | 1130478  | 42.057 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.681 | 200  | 182681   | 33.679 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.428 | 188  | 2092789  | 50.218 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 2568823  | 50.518 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.710 | 264  | 2711852  | 48.124 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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