

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052522\  
 Data File : BP010558.D  
 Acq On : 26 May 2022 03:43  
 Operator : CG/JU  
 Sample : N2883-09  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW4K9

Quant Time: May 26 05:40:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 24 05:50:40 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.881  | 152  | 126743   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.681 | 136  | 555494   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.510 | 164  | 385445   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10      | 17.263 | 188  | 848051   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12          | 21.345 | 240  | 786412   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12          | 23.762 | 264  | 694074   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.334  | 96  | 20488   | 6.890  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.752  | 84  | 306811  | 36.340 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.046  | 99  | 409459  | 38.491 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.210  | 67  | 292929  | 40.985 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.410  | 132 | 326939  | 40.165 | ng/u1 | -0.01 |
| 15) 4-Methylphenol-d8         | 8.587  | 113 | 362437  | 40.091 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.040  | 128 | 183179  | 42.608 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.763  | 143 | 199664  | 43.824 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.304 | 165 | 335694  | 38.967 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.816 | 131 | 377277  | 29.454 | ng/u1 | -0.01 |
| 46) Dimethylphthalate-d6      | 13.922 | 166 | 1326280 | 46.703 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.204 | 160 | 1473488 | 44.983 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.716 | 143 | 173482  | 34.321 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.504 | 176 | 1119856 | 45.327 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.627 | 200 | 171453  | 34.569 | ng/u1 | -0.01 |
| 73) Anthracene-d10            | 17.363 | 188 | 1760987 | 45.858 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.598 | 212 | 2119350 | 50.746 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.609 | 264 | 1624214 | 46.603 | ng/u1 | 0.00  |

Target Compounds Qvalue

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

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