

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072821\  
 Data File : BP006387.D  
 Acq On : 28 Jul 2021 13:31  
 Operator : CG/JU  
 Sample : M3107-16  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BGD16

Quant Time: Jul 28 14:21:57 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jul 24 04:29:12 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.781  | 152  | 337160   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8        | 10.569 | 136  | 1470101  | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10      | 14.410 | 164  | 847980   | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10      | 17.163 | 188  | 1622514  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.263 | 240  | 1349177  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.604 | 264  | 1162633  | 20.000 | ng/ul | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.299  | 96  | 25264   | 2.797  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.711  | 84  | 180869  | 6.745  | ng/ul | 0.00  |
| 7) Phenol-d5                  | 6.958  | 99  | 441801  | 13.952 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67  | 312398  | 15.856 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.316  | 132 | 355874  | 15.127 | ng/ul | -0.01 |
| 15) 4-Methylphenol-d8         | 8.487  | 113 | 338797  | 13.605 | ng/ul | -0.01 |
| 21) Nitrobenzene-d5           | 8.940  | 128 | 175439  | 16.027 | ng/ul | -0.01 |
| 24) 2-Nitrophenol-d4          | 9.652  | 143 | 168416  | 16.114 | ng/ul | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165 | 290193  | 14.169 | ng/ul | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.704 | 131 | 444098  | 13.171 | ng/ul | -0.01 |
| 46) Dimethylphthalate-d6      | 13.828 | 166 | 1044811 | 17.312 | ng/ul | -0.01 |
| 49) Acenaphthylene-d8         | 14.104 | 160 | 1181930 | 15.940 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.610 | 143 | 116993  | 9.763  | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.404 | 176 | 862403  | 16.998 | ng/ul | -0.01 |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200 | 79660   | 8.807  | ng/ul | -0.01 |
| 73) Anthracene-d10            | 17.263 | 188 | 1250214 | 17.069 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.504 | 212 | 1158961 | 16.701 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264 | 682212  | 11.547 | ng/ul | -0.01 |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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