

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072821\  
 Data File : BP006403.D  
 Acq On : 28 Jul 2021 22:38  
 Operator : CG/JU  
 Sample : M3107-13  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BGD05

Quant Time: Jul 29 01:08:51 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  | 287289   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8        | 10.563 | 136  | 1268609  | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10      | 14.404 | 164  | 753453   | 20.000 | ng/u1 | -0.02    |
| 64) Phenanthrene-d10      | 17.157 | 188  | 1477974  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12          | 21.257 | 240  | 1390602  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.598 | 264  | 1322167  | 20.000 | ng/u1 | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.299  | 96  | 21800   | 2.832  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.705  | 84  | 251414  | 11.004 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.952  | 99  | 401471  | 14.879 | ng/u1 | -0.01 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67  | 277919  | 16.554 | ng/u1 | -0.01 |
| 11) 2-Chlorophenol-d4         | 7.316  | 132 | 317447  | 15.836 | ng/u1 | -0.01 |
| 15) 4-Methylphenol-d8         | 8.487  | 113 | 330285  | 15.566 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 8.934  | 128 | 158346  | 16.763 | ng/u1 | -0.02 |
| 24) 2-Nitrophenol-d4          | 9.652  | 143 | 152653  | 16.925 | ng/u1 | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165 | 264143  | 14.945 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.705 | 131 | 419159  | 14.405 | ng/u1 | -0.01 |
| 46) Dimethylphthalate-d6      | 13.828 | 166 | 910468  | 16.978 | ng/u1 | -0.01 |
| 49) Acenaphthylene-d8         | 14.099 | 160 | 1050935 | 15.952 | ng/u1 | -0.01 |
| 54) 4-Nitrophenol-d4          | 14.604 | 143 | 122781  | 11.531 | ng/u1 | -0.01 |
| 60) Fluorene-d10              | 15.398 | 176 | 743207  | 16.486 | ng/u1 | -0.02 |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200 | 83283   | 10.108 | ng/u1 | -0.01 |
| 73) Anthracene-d10            | 17.257 | 188 | 1108073 | 16.608 | ng/u1 | -0.01 |
| 81) Pyrene-d10                | 19.498 | 212 | 1208452 | 16.895 | ng/u1 | -0.01 |
| 92) Benzo(a)pyrene-d12        | 23.445 | 264 | 999586  | 14.877 | ng/u1 | -0.02 |

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

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

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