

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090921\  
 Data File : BP006914.D  
 Acq On : 09 Sep 2021 23:15  
 Operator : CG/JU  
 Sample : M3608-15  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBKL1

Quant Time: Sep 10 01:08:05 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP090821.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 09 11:07:24 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.940  | 152  | 232177   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.740 | 136  | 947205   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.563 | 164  | 581122   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.310 | 188  | 1203583  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12          | 21.386 | 240  | 1211291  | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12          | 23.827 | 264  | 1278846  | 20.000 | ng/ul | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.370  | 96  | 19668   | 3.371  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.793  | 84  | 151994  | 9.997  | ng/ul | 0.01  |
| 7) Phenol-d5                  | 7.093  | 99  | 175035  | 10.257 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.264  | 67  | 465206  | 47.468 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.464  | 132 | 494078  | 35.865 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8         | 8.634  | 113 | 307137  | 22.920 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5           | 9.099  | 128 | 278402  | 48.525 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.816  | 143 | 248934  | 44.071 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.352 | 165 | 507424  | 38.098 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.869 | 131 | 688088  | 37.359 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6      | 13.969 | 166 | 1805355 | 47.966 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8         | 14.257 | 160 | 2199216 | 46.135 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.740 | 143 | 42246   | 6.461  | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.551 | 176 | 1643296 | 49.470 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.663 | 200 | 206203  | 38.728 | ng/ul | 0.00  |
| 73) Anthracene-d10            | 17.410 | 188 | 2793715 | 56.067 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.639 | 212 | 3294891 | 58.082 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.669 | 264 | 3354962 | 54.380 | ng/ul | -0.01 |

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

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

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