

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN062821\  
 Data File : BN015167.D  
 Acq On : 28 Jun 2021 09:58  
 Operator : CG/JU  
 Sample : PB137314BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK314

Quant Time: Jun 28 10:52:57 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061821MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 24 14:43:13 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.716  | 152  | 181711   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8            | 10.493 | 136  | 723803   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.345 | 164  | 418412   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.092 | 188  | 831793   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 824969   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.551 | 264  | 813633   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 31898    | 6.830  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.670  | 84   | 409790   | 32.549 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 508617   | 32.249 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.058  | 67   | 303817   | 32.802 | ng/ul | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.258  | 132  | 405536   | 33.911 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.410  | 113  | 383616   | 31.516 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 8.863  | 128  | 193660   | 39.114 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.581  | 143  | 177819   | 40.318 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.110 | 165  | 345135   | 33.155 | ng/ul | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.622 | 131  | 514780   | 31.909 | ng/ul | -0.01    |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 1014295  | 35.303 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.040 | 160  | 1320186  | 35.396 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.540 | 143  | 144868   | 28.907 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.345 | 176  | 881140   | 35.252 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 68399    | 19.910 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.192 | 188  | 1383040  | 36.328 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.498 | 212  | 1454955  | 33.648 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.404 | 264  | 1489298  | 37.078 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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