

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM111416\  
 Data File : BM007851.D  
 Acq On : 14 Nov 2016 12:54  
 Operator : UM/SJ  
 Sample : PB94659BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK59

Quant Time: Nov 15 02:31:07 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM102016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 15 02:14:40 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.73  | 152  | 158976   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.52 | 136  | 604464   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.37 | 164  | 384539   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.12 | 188  | 749174   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.31 | 240  | 911904   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.56 | 264  | 930526   | 20.00 | ng/ul | 0.00     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 3) 1,4-Dioxane-d8              | 3.28  | 96   | 28310    | 7.63  | ng/uL | 0.00     |
| 5) Phenol-d5                   | 6.91  | 99   | 479587   | 39.09 | ng/ul | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.07  | 67   | 293194   | 40.27 | ng/ul | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.27  | 132  | 382340   | 40.63 | ng/ul | 0.00     |
| 13) 4-Methylphenol-d8          | 8.44  | 113  | 376013   | 39.16 | ng/ul | 0.00     |
| 19) Nitrobenzene-d5            | 8.89  | 128  | 182643   | 41.99 | ng/ul | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.61  | 143  | 200363   | 43.45 | ng/ul | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.15 | 165  | 359278   | 39.63 | ng/ul | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.66 | 131  | 467777   | 45.23 | ng/ul | 0.00     |
| 43) Dimethylphthalate-d6       | 13.79 | 166  | 1100465  | 42.03 | ng/ul | 0.00     |
| 46) Acenaphthylene-d8          | 14.06 | 160  | 1292220  | 41.01 | ng/ul | 0.00     |
| 51) 4-Nitrophenol-d4           | 14.59 | 143  | 139902   | 33.79 | ng/ul | 0.00     |
| 57) Fluorene-d10               | 15.37 | 176  | 935194   | 40.75 | ng/ul | 0.00     |
| 62) 4,6-Dinitro-2-methylphenol | 15.50 | 200  | 126171   | 28.35 | ng/ul | 0.00     |
| 70) Anthracene-d10             | 17.22 | 188  | 1449143  | 42.92 | ng/ul | 0.00     |
| 76) Pyrene-d10                 | 19.52 | 212  | 1669455  | 44.56 | ng/ul | 0.00     |
| 87) Benzo(a)pyrene-d12         | 23.41 | 264  | 1781239  | 43.85 | ng/ul | 0.00     |

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

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

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM111416\  
Data File : BM007851.D  
Acq On : 14 Nov 2016 12:54  
Operator : UM/SJ  
Sample : PB94659BL  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SBLK59

Quant Time: Nov 15 02:31:07 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM102016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Nov 15 02:14:40 2016  
Response via : Initial Calibration

