

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101916\  
 Data File : BM007685.D  
 Acq On : 19 Oct 2016 23:01  
 Operator : UM/SJ  
 Sample : MDL-04-S-ML  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 MDL-04-S-ML

Quant Time: Oct 20 03:13:02 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM101916-MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 20 03:12:33 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 195336   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.53 | 136  | 748379   | 20.00 | ng/ul | 0.00     |
| 13) Acenaphthene-d10      | 14.39 | 164  | 478641   | 20.00 | ng/ul | 0.00     |
| 18) Phenanthrene-d10      | 17.13 | 188  | 972979   | 20.00 | ng/ul | 0.00     |
| 23) Chrysene-d12          | 21.32 | 240  | 1279839  | 20.00 | ng/ul | 0.00     |
| 25) Perylene-d12          | 23.57 | 264  | 1301910  | 20.00 | ng/ul | 0.00     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 2) 1,4-Dioxane-d8              | 3.29  | 96   | 33064    | 7.22  | ng/uL | 0.00     |
| 3) Phenol-d5                   | 6.93  | 99   | 492703   | 32.55 | ng/ul | 0.00     |
| 4) Bis-(2-Chloroethyl)ether-d  | 7.09  | 67   | 300407   | 33.49 | ng/ul | 0.00     |
| 5) 2-Chlorophenol-d4           | 7.29  | 132  | 389915   | 33.56 | ng/ul | 0.00     |
| 6) 4-Methylphenol-d8           | 8.46  | 113  | 382416   | 32.15 | ng/ul | 0.00     |
| 8) Nitrobenzene-d5             | 8.91  | 128  | 178829   | 33.83 | ng/ul | 0.00     |
| 9) 2-Nitrophenol-d4            | 9.63  | 143  | 201161   | 34.86 | ng/ul | 0.00     |
| 10) 2,4-Dichlorophenol-d3      | 10.16 | 165  | 367869   | 32.40 | ng/ul | 0.00     |
| 11) 4-Chloroaniline-d4         | 10.68 | 131  | 465884   | 36.16 | ng/ul | 0.00     |
| 14) Dimethylphthalate-d6       | 13.80 | 166  | 1120002  | 33.54 | ng/ul | 0.00     |
| 15) Acenaphthylene-d8          | 14.08 | 160  | 1330374  | 33.99 | ng/ul | 0.00     |
| 16) 4-Nitrophenol-d4           | 14.60 | 143  | 154319   | 28.11 | ng/ul | 0.00     |
| 17) Fluorene-d10               | 15.38 | 176  | 965834   | 33.52 | ng/ul | 0.00     |
| 19) 4,6-Dinitro-2-methylphenol | 15.51 | 200  | 167718   | 28.45 | ng/ul | 0.00     |
| 20) Anthracene-d10             | 17.23 | 188  | 1453571  | 32.86 | ng/ul | 0.00     |
| 24) Pyrene-d10                 | 19.53 | 212  | 1732512  | 33.44 | ng/ul | 0.00     |
| 26) Benzo(a)pyrene-d12         | 23.43 | 264  | 1917004  | 33.81 | ng/ul | 0.00     |

| Target Compounds               | R.T.  | QIon | Response | Conc | Units | Qvalue |
|--------------------------------|-------|------|----------|------|-------|--------|
| 12) 1-Methylnaphthalene        | 12.42 | 142  | 79499    | 3.09 | ng/ul | 94     |
| 21) 1,2,3,4-Tetrachlorobenzene | 13.19 | 216  | 44195    | 3.12 | ng/uL | 98     |
| 22) Pentachlorobenzene         | 14.71 | 250  | 47572    | 3.21 | ng/uL | 95     |

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

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