

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101916\  
 Data File : BM007695.D  
 Acq On : 20 Oct 2016 06:00  
 Operator : UM/SJ  
 Sample : MDL-05-W  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 MDL-05-W

Manual Integrations  
 APPROVED

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 10/20/2016 5:04:35 PM

Quant Time: Oct 20 07:10:30 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM101916-MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 20 03:26:07 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 245760   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.53 | 136  | 951979   | 20.00 | ng/ul | 0.00     |
| 13) Acenaphthene-d10      | 14.39 | 164  | 604191   | 20.00 | ng/ul | 0.00     |
| 18) Phenanthrene-d10      | 17.13 | 188  | 1203002m | 20.00 | ng/ul | 0.00     |
| 23) Chrysene-d12          | 21.32 | 240  | 1538258  | 20.00 | ng/ul | 0.00     |
| 25) Perylene-d12          | 23.57 | 264  | 1566972  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.29  | 96  | 39100   | 6.79  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.93  | 99  | 580450  | 30.48 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 7.09  | 67  | 356440  | 31.58 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.29  | 132 | 459837  | 31.46 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.46  | 113 | 450275  | 30.09 | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.91  | 128 | 216003  | 32.12 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.63  | 143 | 240483  | 32.76 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.16 | 165 | 432733  | 29.96 | ng/ul | 0.00 |
| 11) 4-Chloroaniline-d4         | 10.67 | 131 | 550027  | 33.56 | ng/ul | 0.00 |
| 14) Dimethylphthalate-d6       | 13.80 | 166 | 1296899 | 30.77 | ng/ul | 0.00 |
| 15) Acenaphthylene-d8          | 14.08 | 160 | 1557203 | 31.52 | ng/ul | 0.00 |
| 16) 4-Nitrophenol-d4           | 14.60 | 143 | 187131  | 27.00 | ng/ul | 0.00 |
| 17) Fluorene-d10               | 15.38 | 176 | 1126389 | 30.97 | ng/ul | 0.00 |
| 19) 4,6-Dinitro-2-methylphenol | 15.51 | 200 | 197237  | 27.06 | ng/ul | 0.00 |
| 20) Anthracene-d10             | 17.23 | 188 | 1677882 | 30.68 | ng/ul | 0.00 |
| 24) Pyrene-d10                 | 19.53 | 212 | 1968302 | 31.61 | ng/ul | 0.00 |
| 26) Benzo(a)pyrene-d12         | 23.43 | 264 | 2129002 | 31.19 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |      |       | Qvalue |
|--------------------------------|-------|-----|--------|------|-------|--------|
| 12) 1-Methylnaphthalene        | 12.42 | 142 | 94901m | 2.90 | ng/ul |        |
| 21) 1,2,3,4-Tetrachlorobenzene | 13.18 | 216 | 52688  | 3.01 | ng/uL | 97     |
| 22) Pentachlorobenzene         | 14.71 | 250 | 55116  | 3.01 | ng/uL | 99     |

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

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