

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM073116\  
 Data File : BM006811.D  
 Acq On : 01 Aug 2016 04:59  
 Operator : UM/SJ  
 Sample : H4189-15  
 Misc :  
 ALS Vial : 79 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9ZQ4

Manual Integrations  
 APPROVED

sohil  
 8/1/2016 7:08:40 PM

Quant Time: Aug 01 08:13:25 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jul 30 01:37:26 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.60  | 152  | 227256   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.38 | 136  | 994034   | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.26 | 164  | 551432   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.02 | 188  | 1228696  | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.22 | 240  | 1301325  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.42 | 264  | 1371782  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.13  | 96  | 6951    | 1.24  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.79  | 99  | 292628  | 15.14 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 206413  | 17.23 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.14  | 132 | 247523  | 16.02 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.32  | 113 | 168292  | 11.22 | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.76  | 128 | 130845  | 17.39 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.48  | 143 | 146544  | 18.31 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 245628  | 16.24 | ng/ul | 0.00 |
| 12) 4-Chloroaniline-d4         | 10.53 | 131 | 264752  | 15.90 | ng/ul | 0.00 |
| 16) Dimethylphthalate-d6       | 13.67 | 166 | 776576  | 17.96 | ng/ul | 0.00 |
| 17) Acenaphthylene-d8          | 13.95 | 160 | 1007479 | 18.26 | ng/ul | 0.00 |
| 20) 4-Nitrophenol-d4           | 14.50 | 143 | 98361   | 13.10 | ng/ul | 0.01 |
| 21) Fluorene-d10               | 15.26 | 176 | 695891  | 18.00 | ng/ul | 0.00 |
| 24) 4,6-Dinitro-2-methylphenol | 15.41 | 200 | 75915   | 11.90 | ng/ul | 0.01 |
| 26) Anthracene-d10             | 17.11 | 188 | 1052508 | 18.04 | ng/ul | 0.00 |
| 30) Pyrene-d10                 | 19.42 | 212 | 1165244 | 19.65 | ng/ul | 0.00 |
| 37) Benzo(a)pyrene-d12         | 23.29 | 264 | 1138112 | 18.04 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        |      |       | Qvalue |
|----------------------------|-------|-----|--------|------|-------|--------|
| 25) Phenanthrene           | 17.06 | 178 | 144788 | 2.10 | ng/ul | 98     |
| 28) Fluoranthene           | 19.08 | 202 | 365968 | 4.51 | ng/ul | 99     |
| 31) Pyrene                 | 19.44 | 202 | 324511 | 4.12 | ng/ul | 99     |
| 32) Benzo(a)anthracene     | 21.20 | 228 | 188713 | 2.39 | ng/ul | 95     |
| 33) Chrysene               | 21.26 | 228 | 206730 | 2.84 | ng/ul | 95     |
| 35) Benzo(b)fluoranthene   | 22.77 | 252 | 290753 | 3.54 | ng/ul | 96     |
| 36) Benzo(k)fluoranthene   | 22.80 | 252 | 97583m | 1.23 | ng/ul |        |
| 38) Benzo(a)pyrene         | 23.33 | 252 | 203510 | 2.54 | ng/ul | 96     |
| 39) Indeno(1,2,3-cd)pyrene | 25.63 | 276 | 156546 | 1.69 | ng/ul | 98     |
| 41) Benzo(g,h,i)perylene   | 26.31 | 276 | 152011 | 1.91 | ng/ul | 97     |

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

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