

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM073116\  
 Data File : BM006806.D  
 Acq On : 01 Aug 2016 01:56  
 Operator : UM/SJ  
 Sample : H4188-15MS  
 Misc :  
 ALS Vial : 74 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9ZP2MS

Manual Integrations  
 APPROVED

sohil  
 8/1/2016 7:08:00 PM

Quant Time: Aug 01 08:00:18 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  | 255591   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.38 | 136  | 1034095  | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.26 | 164  | 558143   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.01 | 188  | 1238490  | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.22 | 240  | 1340316  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.42 | 264  | 1426690  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.13  | 96  | 10078   | 1.59  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.79  | 99  | 318515  | 14.65 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 229594  | 17.04 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.14  | 132 | 277974  | 16.00 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.32  | 113 | 160404  | 9.51  | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.76  | 128 | 144547  | 18.47 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.47  | 143 | 158381  | 19.03 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 271099  | 17.23 | ng/ul | 0.00 |
| 12) 4-Chloroaniline-d4         | 10.53 | 131 | 232928  | 13.45 | ng/ul | 0.00 |
| 16) Dimethylphthalate-d6       | 13.67 | 166 | 839121  | 19.17 | ng/ul | 0.00 |
| 17) Acenaphthylene-d8          | 13.95 | 160 | 1085630 | 19.44 | ng/ul | 0.00 |
| 20) 4-Nitrophenol-d4           | 14.49 | 143 | 81979   | 10.79 | ng/ul | 0.00 |
| 21) Fluorene-d10               | 15.26 | 176 | 754776  | 19.28 | ng/ul | 0.00 |
| 24) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 73802   | 11.48 | ng/ul | 0.00 |
| 26) Anthracene-d10             | 17.11 | 188 | 1154253 | 19.63 | ng/ul | 0.00 |
| 30) Pyrene-d10                 | 19.42 | 212 | 1279252 | 20.95 | ng/ul | 0.00 |
| 37) Benzo(a)pyrene-d12         | 23.28 | 264 | 1290508 | 19.66 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |       |        | Qvalue |
|----------------------------|-------|-----|---------|-------|--------|--------|
| 19) Acenaphthene           | 14.32 | 153 | 630385  | 16.40 | ng/ul  | 94     |
| 25) Phenanthrene           | 17.05 | 178 | 203796  | 2.93  | ng/ul  | 99     |
| 28) Fluoranthene           | 19.08 | 202 | 488574  | 5.97  | ng/ul  | 97     |
| 31) Pyrene                 | 19.44 | 202 | 1788531 | 22.04 | ng/ul  | 97     |
| 32) Benzo(a)anthracene     | 21.20 | 228 | 267611  | 3.29  | ng/ul  | 97     |
| 33) Chrysene               | 21.25 | 228 | 275672m | 3.68  | ng/ul  |        |
| 35) Benzo(b)fluoranthene   | 22.76 | 252 | 407895  | 4.77  | ng/ul  | 98     |
| 36) Benzo(k)fluoranthene   | 22.79 | 252 | 152886m | 1.85  | ng/ul  |        |
| 38) Benzo(a)pyrene         | 23.32 | 252 | 284137  | 3.41  | ng/ul  | 96     |
| 39) Indeno(1,2,3-cd)pyrene | 25.63 | 276 | 229621  | 2.38  | ng/ul  | 98     |
| 41) Benzo(g,h,i)perylene   | 26.30 | 276 | 228007  | 2.76  | ng/ul# | 95     |

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

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