

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM062616\  
 Data File : BM006036.D  
 Acq On : 26 Jun 2016 12:36  
 Operator : UM/SJ  
 Sample : H3669-14  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9Y67

Manual Integrations  
 APPROVED

sohil  
 6/27/2016 8:05:50 PM

Quant Time: Jun 27 05:28:10 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM062516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 27 05:15:31 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.66  | 152  | 203582   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.43 | 136  | 1022492  | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.30 | 164  | 671443   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.05 | 188  | 1683971  | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.24 | 240  | 1959242  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.46 | 264  | 1972373  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.18  | 96  | 3530    | 0.77  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.83  | 99  | 216345  | 10.10 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.99  | 67  | 132905  | 11.15 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.19  | 132 | 160004  | 10.73 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.36  | 113 | 158137  | 8.75  | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.80  | 128 | 84590   | 10.87 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.52  | 143 | 90392   | 10.44 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.06 | 165 | 189898  | 11.43 | ng/ul | 0.00 |
| 12) 4-Chloroaniline-d4         | 10.57 | 131 | 156706  | 8.69  | ng/ul | 0.00 |
| 16) Dimethylphthalate-d6       | 13.72 | 166 | 657022  | 11.43 | ng/ul | 0.00 |
| 17) Acenaphthylene-d8          | 13.99 | 160 | 824138  | 12.36 | ng/ul | 0.00 |
| 20) 4-Nitrophenol-d4           | 14.50 | 143 | 102632  | 8.37  | ng/ul | 0.00 |
| 21) Fluorene-d10               | 15.30 | 176 | 618430  | 12.60 | ng/ul | 0.00 |
| 24) 4,6-Dinitro-2-methylphenol | 15.42 | 200 | 73041   | 7.81  | ng/ul | 0.00 |
| 26) Anthracene-d10             | 17.15 | 188 | 1002888 | 12.97 | ng/ul | 0.00 |
| 30) Pyrene-d10                 | 19.44 | 212 | 1202073 | 14.01 | ng/ul | 0.00 |
| 37) Benzo(a)pyrene-d12         | 23.32 | 264 | 1160774 | 13.04 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |       |        | Qvalue |
|----------------------------|-------|-----|---------|-------|--------|--------|
| 25) Phenanthrene           | 17.09 | 178 | 682474  | 7.53  | ng/ul  | 99     |
| 27) Anthracene             | 17.18 | 178 | 110404  | 1.18  | ng/ul  | 100    |
| 28) Fluoranthene           | 19.11 | 202 | 1195984 | 10.31 | ng/ul  | 98     |
| 31) Pyrene                 | 19.47 | 202 | 954327  | 8.51  | ng/ul  | 98     |
| 32) Benzo(a)anthracene     | 21.23 | 228 | 533955  | 4.59  | ng/ul  | 98     |
| 33) Chrysene               | 21.28 | 228 | 587419  | 5.52  | ng/ul  | 96     |
| 35) Benzo(b)fluoranthene   | 22.80 | 252 | 807704  | 7.05  | ng/ul# | 96     |
| 36) Benzo(k)fluoranthene   | 22.84 | 252 | 270868m | 2.54  | ng/ul  |        |
| 38) Benzo(a)pyrene         | 23.36 | 252 | 509074  | 4.63  | ng/ul  | 97     |
| 39) Indeno(1,2,3-cd)pyrene | 25.68 | 276 | 370199  | 2.83  | ng/ul  | 98     |
| 41) Benzo(g,h,i)perylene   | 26.37 | 276 | 338567  | 2.99  | ng/ul  | 99     |

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

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