

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011616\  
 Data File : BM004052.D  
 Acq On : 15 Jan 2016 17:48  
 Operator : SJ/UM  
 Sample : H1100-11  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BC992

Manual Integrations  
 APPROVED

Sohil  
 1/16/2016 5:01:16 AM

Quant Time: Jan 16 02:50:01 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jan 16 00:22:06 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.66  | 152  | 39878    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.45 | 136  | 191840   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.32 | 164  | 110287   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.07 | 188  | 242416   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.28 | 240  | 207057   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.52 | 264  | 183529   | 20.00 | ng/ul | 0.01     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.13  | 96  | 2805   | 3.15  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.85  | 99  | 68028  | 20.55 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.00  | 67  | 40637  | 21.71 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.20  | 132 | 59273  | 21.78 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.38  | 113 | 41758  | 15.56 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.82  | 128 | 31107  | 21.53 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.55  | 143 | 34249  | 21.62 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.08 | 165 | 59369  | 21.09 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.60 | 131 | 62875  | 16.92 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.73 | 166 | 200713 | 23.18 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.01 | 160 | 248012 | 23.51 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.55 | 143 | 25892  | 16.65 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.32 | 176 | 177595 | 24.18 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.46 | 200 | 14609  | 10.69 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.17 | 188 | 273161 | 24.36 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.47 | 212 | 285280 | 26.60 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.37 | 264 | 230265 | 25.12 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        |       |        | Qvalue |
|----------------------------|-------|-----|--------|-------|--------|--------|
| 45) Dimethylphthalate      | 13.78 | 163 | 87105  | 9.87  | ng/ul  | 100    |
| 70) Phenanthrene           | 17.12 | 178 | 140010 | 10.27 | ng/ul  | 99     |
| 72) Anthracene             | 17.20 | 178 | 35684  | 2.55  | ng/ul  | 99     |
| 77) Fluoranthene           | 19.14 | 202 | 288439 | 20.36 | ng/ul  | 99     |
| 80) Pyrene                 | 19.51 | 202 | 325562 | 23.16 | ng/ul  | 100    |
| 83) Benzo(a)anthracene     | 21.26 | 228 | 138747 | 11.25 | ng/ul  | 98     |
| 85) Chrysene               | 21.32 | 228 | 149229 | 12.74 | ng/ul  | 98     |
| 88) Benzo(b)fluoranthene   | 22.85 | 252 | 156197 | 14.00 | ng/ul  | 98     |
| 89) Benzo(k)fluoranthene   | 22.87 | 252 | 54458m | 5.13  | ng/ul  |        |
| 91) Benzo(a)pyrene         | 23.42 | 252 | 119813 | 11.32 | ng/ul  | 97     |
| 92) Indeno(1,2,3-cd)pyrene | 25.77 | 276 | 73243  | 6.26  | ng/ul  | 96     |
| 93) Dibenzo(a,h)anthracene | 25.77 | 278 | 21162  | 2.15  | ng/ul# | 86     |
| 94) Benzo(g,h,i)perylene   | 26.46 | 276 | 63835  | 6.49  | ng/ul  | 97     |

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

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