

Data Path : Z:\HPCHEM1\BNA M\DATA\BM080416\  
 Data File : BM006890.D  
 Acq On : 05 Aug 2016 07:14  
 Operator : UM/SJ  
 Sample : H4224-09  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BDJL9

Manual Integrations  
 APPROVED

umangi  
 8/5/2016 7:03:18 PM

Quant Time: Aug 05 07:51:59 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM080316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 05 02:18:21 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 284803   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.33 | 136  | 1185877  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.22 | 164  | 691325   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.97 | 188  | 1640700  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.19 | 240  | 1913658  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.37 | 264  | 1853369  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.11  | 96  | 8781    | 1.70  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.79  | 99  | 310749  | 14.82 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.91  | 67  | 217457  | 15.53 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.12  | 132 | 270939  | 16.36 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.32  | 113 | 244998  | 14.37 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.73  | 128 | 120187  | 14.82 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 137810  | 14.28 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.00 | 165 | 247710m | 12.41 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.52 | 131 | 316732m | 15.36 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.64 | 166 | 966222  | 16.97 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.91 | 160 | 1118635 | 16.08 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.55 | 143 | 84272m  | 6.75  | ng/ul | 0.01 |
| 57) Fluorene-d10               | 15.22 | 176 | 849332  | 16.07 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.39 | 200 | 82558   | 7.71  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.07 | 188 | 1357822 | 17.23 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.38 | 212 | 1681964 | 17.18 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.23 | 264 | 1597135 | 17.70 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |      |        | Ovalue |
|--------------------------------|-------|-----|--------|------|--------|--------|
| 44) Dimethylphthalate          | 13.69 | 163 | 433932 | 7.67 | ng/ul  | 98     |
| 69) Phenanthrene               | 17.02 | 178 | 177601 | 1.96 | ng/ul# | 95     |
| 74) Fluoranthene               | 19.05 | 202 | 389834 | 3.36 | ng/ul# | 96     |
| 77) Pyrene                     | 19.41 | 202 | 367708 | 2.92 | ng/ul# | 94     |
| 80) Benzo(a)anthracene         | 21.17 | 228 | 183362 | 1.57 | ng/ul  | 99     |
| 81) Bis(2-ethylhexyl)phthalate | 21.09 | 149 | 75114  | 1.26 | ng/ul  | 99     |
| 82) Chrysene                   | 21.22 | 228 | 181474 | 1.77 | ng/ul  | 97     |
| 85) Benzo(b)fluoranthene       | 22.73 | 252 | 197625 | 1.69 | ng/ul# | 91     |
| 88) Benzo(a)pyrene             | 23.27 | 252 | 151196 | 1.36 | ng/ul# | 94     |

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

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