

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022616\  
 Data File : BM004433.D  
 Acq On : 26 Feb 2016 18:52  
 Operator : SJ/UM  
 Sample : H1564-13MS  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9R26MS

Manual Integrations  
 APPROVED

UMANGI  
 2/29/2016 2:50:34 PM

Quant Time: Feb 27 00:36:44 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM022216.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Feb 26 23:20:20 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.62  | 152  | 22761    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.40 | 136  | 98504    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.27 | 164  | 54609    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.03 | 188  | 113971   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.24 | 240  | 119992   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.46 | 264  | 119611   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0      | 0.00  | ng/uL |      |
| 5) Phenol-d5                   | 6.81  | 99  | 5869   | 3.08  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.96  | 67  | 2870   | 2.96  | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.16  | 132 | 5927   | 3.60  | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.35  | 113 | 3539   | 2.08  | ng/ul | 0.01 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 1671   | 2.23  | ng/ul | 0.01 |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 1618   | 1.89  | ng/ul | 0.01 |
| 26) 2,4-Dichlorophenol-d3      | 10.06 | 165 | 3387m  | 2.19  | ng/ul | 0.03 |
| 29) 4-Chloroaniline-d4         | 10.57 | 131 | 3176m  | 1.65  | ng/ul | 0.03 |
| 44) Dimethylphthalate-d6       | 13.69 | 166 | 15793  | 3.37  | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.97 | 160 | 17611  | 3.21  | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.54 | 143 | 2317   | 3.08  | ng/ul | 0.02 |
| 58) Fluorene-d10               | 15.28 | 176 | 16179  | 4.06  | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 0.00  | 200 | 0d     | 0.00  | ng/ul |      |
| 71) Anthracene-d10             | 17.13 | 188 | 32529  | 5.82  | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.44 | 212 | 76023  | 13.10 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.32 | 264 | 104944 | 16.72 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 6) Phenol                      | 6.84  | 94  | 5624   | 2.76  | ng/ul  | 95     |
| 10) 2-Chlorophenol             | 7.19  | 128 | 5335   | 3.12  | ng/ul  | 97     |
| 15) N-Nitroso-di-n-propylamine | 8.42  | 70  | 2147   | 1.67  | ng/ul# | 90     |
| 33) 4-Chloro-3-methylphenol    | 11.74 | 107 | 2725   | 1.43  | ng/ul  | 92     |
| 45) Dimethylphthalate          | 13.74 | 163 | 5198   | 1.06  | ng/ul# | 95     |
| 50) Acenaphthene               | 14.34 | 153 | 12431  | 3.05  | ng/ul  | 98     |
| 53) 4-Nitrophenol              | 14.56 | 109 | 1261   | 2.24  | ng/ul  | 93     |
| 55) 2,4-Dinitrotoluene         | 14.68 | 165 | 3425   | 2.34  | ng/ul# | 94     |
| 69) Pentachlorophenol          | 16.70 | 266 | 1358   | 2.49  | ng/ul  | 97     |
| 77) Fluoranthene               | 19.10 | 202 | 17685  | 2.28  | ng/ul  | 99     |
| 80) Pyrene                     | 19.47 | 202 | 105377 | 13.44 | ng/ul  | 100    |
| 83) Benzo(a)anthracene         | 21.23 | 228 | 7729   | 1.00  | ng/ul  | 94     |
| 85) Chrysene                   | 21.28 | 228 | 9643   | 1.30  | ng/ul  | 98     |
| 88) Benzo(b)fluoranthene       | 22.80 | 252 | 12465  | 1.54  | ng/ul# | 94     |
| 91) Benzo(a)pyrene             | 23.37 | 252 | 8565   | 1.14  | ng/ul  | 97     |

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

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