

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM010617\  
 Data File : BM008735.D  
 Acq On : 06 Jan 2017 15:02  
 Operator : UM/SJ  
 Sample : SSTD16055  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD16055

Manual Integrations  
 APPROVED

Sohil  
 1/9/2017 10:45:04 AM

Quant Time: Jan 06 16:06:15 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 06 13:55:44 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.96  | 152  | 114187   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.77 | 136  | 497761   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.60 | 164  | 389225   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.34 | 188  | 879973   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.51 | 240  | 1045701  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.88 | 264  | 963915   | 20.00 | ng/ul | 0.01     |

## System Monitoring Compounds

|                                |       |     |         |        |       |      |
|--------------------------------|-------|-----|---------|--------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d      | 0.00   | ng/uL |      |
| 5) Phenol-d5                   | 7.13  | 99  | 1571819 | 184.64 | ng/ul | 0.02 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.30  | 67  | 1027630 | 209.98 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 0.00  | 132 | 0d      | 0.00   | ng/ul |      |
| 13) 4-Methylphenol-d8          | 8.67  | 113 | 1211732 | 183.15 | ng/ul | 0.02 |
| 19) Nitrobenzene-d5            | 0.00  | 128 | 0d      | 0.00   | ng/ul |      |
| 22) 2-Nitrophenol-d4           | 0.00  | 143 | 0d      | 0.00   | ng/ul |      |
| 26) 2,4-Dichlorophenol-d3      | 0.00  | 165 | 0d      | 0.00   | ng/ul |      |
| 29) 4-Chloroaniline-d4         | 10.91 | 131 | 1118623 | 155.14 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 0.00  | 166 | 0d      | 0.00   | ng/ul |      |
| 46) Acenaphthylene-d8          | 0.00  | 160 | 0d      | 0.00   | ng/ul |      |
| 51) 4-Nitrophenol-d4           | 14.80 | 143 | 738667  | 190.60 | ng/ul | 0.03 |
| 57) Fluorene-d10               | 0.00  | 176 | 0d      | 0.00   | ng/ul |      |
| 62) 4,6-Dinitro-2-methylphenol | 15.72 | 200 | 868495  | 166.22 | ng/ul | 0.02 |
| 70) Anthracene-d10             | 0.00  | 188 | 0d      | 0.00   | ng/ul |      |
| 76) Pyrene-d10                 | 0.00  | 212 | 0d      | 0.00   | ng/ul |      |
| 87) Benzo(a)pyrene-d12         | 0.00  | 264 | 0d      | 0.00   | ng/ul |      |

## Target Compounds

|                                |       |     |         |        |        | Qvalue |
|--------------------------------|-------|-----|---------|--------|--------|--------|
| 4) Benzaldehyde                | 7.11  | 77  | 626049m | 101.62 | ng/ul  |        |
| 6) Phenol                      | 7.15  | 94  | 1691593 | 193.54 | ng/ul# | 73     |
| 8) Bis(2-Chloroethyl)ether     | 7.39  | 93  | 1213311 | 182.92 | ng/ul# | 88     |
| 11) 2-Methylphenol             | 8.40  | 108 | 1246475 | 192.35 | ng/ul  | 94     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.50  | 45  | 1810394 | 245.56 | ng/ul# | 93     |
| 14) Acetophenone               | 8.80  | 105 | 2110449 | 199.41 | ng/ul# | 85     |
| 16) 4-Methylphenol             | 8.75  | 108 | 1319781 | 189.01 | ng/ul  | 97     |
| 30) 4-Chloroaniline            | 10.93 | 127 | 1159317 | 156.09 | ng/ul  | 99     |
| 32) Caprolactam                | 11.77 | 113 | 394614m | 195.36 | ng/ul  |        |
| 37) Hexachlorocyclopentadiene  | 12.77 | 237 | 1408587 | 235.37 | ng/ul  | 97     |
| 48) 3-Nitroaniline             | 14.52 | 138 | 707773  | 154.72 | ng/ul  | 87     |
| 50) 2,4-Dinitrophenol          | 14.71 | 184 | 609823  | 198.00 | ng/ul# | 66     |
| 52) 4-Nitrophenol              | 14.82 | 109 | 1201310 | 329.73 | ng/ul# | 70     |
| 60) 4-Nitroaniline             | 15.69 | 138 | 835104  | 182.05 | ng/ul# | 41     |
| 63) 4,6-Dinitro-2-methylphenol | 15.73 | 198 | 892781  | 166.14 | ng/ul# | 79     |
| 67) Atrazine                   | 16.81 | 200 | 1617820 | 168.11 | ng/ul  | 96     |
| 68) Pentachlorophenol          | 16.99 | 266 | 1125475 | 180.20 | ng/ul  | 98     |
| 72) Carbazole                  | 17.75 | 167 | 6390906 | 163.90 | ng/ul  | 98     |
| 74) Fluoranthene               | 19.39 | 202 | 9053747 | 177.19 | ng/ul  | 96     |
| 79) 3,3'-Dichlorobenzidine     | 21.43 | 252 | 2979106 | 161.37 | ng/ul  | 100    |
| 84) Di-n-octyl phthalate       | 22.33 | 149 | 8948764 | 180.33 | ng/ul# | 94     |

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

