

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040116\  
 Data File : BM004843.D  
 Acq On : 01 Apr 2016 16:12  
 Operator : UM/SJ  
 Sample : SSTD00550  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD00551

Manual Integrations  
 APPROVED

Sohil  
 4/4/2016 6:33:59 PM

Quant Time: Apr 01 18:23:31 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM040116.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 01 18:03:21 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.83  | 152  | 42869    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.62 | 136  | 175979   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.46 | 164  | 121799   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.21 | 188  | 248391   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.41 | 240  | 1010455  | 20.00 | ng/ul | 0.01     |
| 86) Perylene-d12          | 23.71 | 264  | 371858   | 20.00 | ng/ul | 0.02     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.32  | 96  | 2261   | 1.82 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 0.00  | 99  | 0d     | 0.00 | ng/ul |      |
| 7) Bis-(2-Chloroethyl)ether-d  | 0.00  | 67  | 0d     | 0.00 | ng/ul |      |
| 9) 2-Chlorophenol-d4           | 7.36  | 132 | 13336  | 4.70 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 0.00  | 113 | 0d     | 0.00 | ng/ul |      |
| 19) Nitrobenzene-d5            | 8.99  | 128 | 6993   | 5.71 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.70  | 143 | 7753   | 5.87 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.25 | 165 | 14222  | 5.48 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d     | 0.00 | ng/ul |      |
| 44) Dimethylphthalate-d6       | 13.87 | 166 | 48285  | 5.07 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.16 | 160 | 59086  | 5.06 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 0.00  | 143 | 0d     | 0.00 | ng/ul |      |
| 58) Fluorene-d10               | 15.46 | 176 | 42369  | 5.11 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 0.00  | 200 | 0d     | 0.00 | ng/ul |      |
| 71) Anthracene-d10             | 17.31 | 188 | 67081m | 5.87 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.61 | 212 | 80672  | 1.75 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.56 | 264 | 85797  | 5.17 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |       |      |        | Ovalue |
|--------------------------------|-------|-----|-------|------|--------|--------|
| 2) 1,4-Dioxane                 | 3.35  | 88  | 2704  | 1.87 | ng/uL# | 69     |
| 10) 2-Chlorophenol             | 7.39  | 128 | 14456 | 4.83 | ng/ul# | 91     |
| 15) N-Nitroso-di-n-propylamine | 8.62  | 70  | 14107 | 6.27 | ng/ul# | 88     |
| 17) Hexachloroethane           | 8.90  | 117 | 5735  | 4.82 | ng/ul  | 99     |
| 20) Nitrobenzene               | 9.03  | 77  | 16394 | 5.43 | ng/ul  | 99     |
| 21) Isophorone                 | 9.55  | 82  | 34873 | 6.03 | ng/ul  | 98     |
| 23) 2-Nitrophenol              | 9.73  | 139 | 8286  | 5.66 | ng/ul# | 89     |
| 24) 2,4-Dimethylphenol         | 9.80  | 107 | 18105 | 5.65 | ng/ul  | 95     |
| 25) Bis(2-Chloroethoxy)methane | 10.03 | 93  | 21539 | 5.77 | ng/ul# | 88     |
| 27) 2,4-Dichlorophenol         | 10.27 | 162 | 14334 | 5.37 | ng/ul  | 96     |
| 28) Naphthalene                | 10.67 | 128 | 47156 | 5.14 | ng/ul  | 100    |
| 31) Hexachlorobutadiene        | 10.95 | 225 | 9056  | 5.68 | ng/ul  | 99     |
| 33) 4-Chloro-3-methylphenol    | 11.90 | 107 | 18931 | 6.26 | ng/ul  | 97     |
| 34) 2-Methylnaphthalene        | 12.28 | 142 | 36512 | 5.59 | ng/ul  | 99     |
| 35) 1-Methylnaphthalene        | 12.50 | 142 | 35602 | 5.67 | ng/ul  | 100    |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.65 | 216 | 18559 | 4.77 | ng/ul  | 99     |
| 39) 2,4,6-Trichlorophenol      | 12.89 | 196 | 11065 | 4.31 | ng/ul  | 99     |
| 40) 2,4,5-Trichlorophenol      | 12.96 | 196 | 13418 | 4.84 | ng/ul  | 97     |
| 41) 1,1'-Biphenyl              | 13.29 | 154 | 52405 | 5.15 | ng/ul  | 98     |
| 42) 2-Chloronaphthalene        | 13.34 | 162 | 36601 | 4.78 | ng/ul  | 96     |
| 43) 2-Nitroaniline             | 13.55 | 65  | 10447 | 5.07 | ng/ul  | 88     |
| 45) Dimethylphthalate          | 13.92 | 163 | 50113 | 5.18 | ng/ul  | 99     |
| 46) 2,6-Dinitrotoluene         | 14.04 | 165 | 9582  | 5.57 | ng/ul# | 79     |

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 Sample : SSTD00550  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD00551

Manual Integrations  
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Quant Time: Apr 01 18:23:31 2016  
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 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 48) Acenaphthylene             | 14.19 | 152  | 61130    | 4.87  | ng/ul  | 100      |
| 50) Acenaphthene               | 14.53 | 153  | 42048    | 4.98  | ng/ul  | 100      |
| 54) Dibenzofuran               | 14.86 | 168  | 60032    | 4.99  | ng/ul  | 98       |
| 55) 2,4-Dinitrotoluene         | 14.83 | 165  | 12219    | 4.71  | ng/ul# | 91       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.09 | 232  | 12884    | 5.28  | ng/ul# | 95       |
| 57) Diethylphthalate           | 15.28 | 149  | 58913    | 5.98  | ng/ul  | 98       |
| 59) Fluorene                   | 15.52 | 166  | 51546    | 5.37  | ng/ul  | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.51 | 204  | 23815    | 5.20  | ng/ul  | 93       |
| 65) N-Nitrosodiphenylamine     | 15.72 | 169  | 44139    | 5.87  | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.40 | 248  | 14898    | 5.96  | ng/ul  | 91       |
| 67) Hexachlorobenzene          | 16.52 | 284  | 15163    | 5.40  | ng/ul  | 97       |
| 70) Phenanthrene               | 17.25 | 178  | 82958    | 5.85  | ng/ul  | 99       |
| 72) Anthracene                 | 17.35 | 178  | 82474    | 5.77  | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.26 | 216  | 18553    | 5.43  | ng/uL  | 96       |
| 74) Pentachlorobenzene         | 14.78 | 250  | 16434    | 4.98  | ng/uL  | 99       |
| 76) Di-n-butylphthalate        | 18.17 | 149  | 145843   | 8.05  | ng/ul  | 98       |
| 80) Pyrene                     | 19.64 | 202  | 100363   | 1.64  | ng/ul# | 95       |
| 81) Butylbenzylphthalate       | 20.54 | 149  | 108333   | 4.42  | ng/ul  | 96       |
| 83) Benzo(a)anthracene         | 21.39 | 228  | 121812   | 2.07  | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.32 | 149  | 813948   | 23.37 | ng/ul# | 96       |
| 85) Chrysene                   | 21.44 | 228  | 745778   | 13.24 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.01 | 252  | 111694   | 5.04  | ng/ul# | 93       |
| 89) Benzo(k)fluoranthene       | 23.06 | 252  | 107878   | 5.12  | ng/ul# | 92       |
| 91) Benzo(a)pyrene             | 23.60 | 252  | 108385   | 5.10  | ng/ul# | 92       |
| 92) Indeno(1,2,3-cd)pyrene     | 26.03 | 276  | 130736   | 5.47  | ng/ul  | 99       |
| 93) Dibenzo(a,h)anthracene     | 26.04 | 278  | 108928   | 5.51  | ng/ul  | 99       |
| 94) Benzo(g,h,i)perylene       | 26.74 | 276  | 110465   | 5.46  | ng/ul  | 97       |

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

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