

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM091018\  
 Data File : BM016710.D  
 Acq On : 10 Sep 2018 14:52  
 Operator : SJ/JU  
 Sample : SSTD16052  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD16052

Quant Time: Sep 10 15:23:19 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 10 13:31:33 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.75  | 152  | 195209   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.53 | 136  | 807844   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.39 | 164  | 427734   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.13 | 188  | 926166   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.33 | 240  | 987924   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.57 | 264  | 1031825  | 20.00 | ng/ul | 0.01     |

## System Monitoring Compounds

|                                |       |     |         |        |       |      |
|--------------------------------|-------|-----|---------|--------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.27  | 96  | 282174  | 65.02  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.92  | 99  | 2639482 | 140.82 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.09  | 67  | 1592712 | 126.09 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.29  | 132 | 2185556 | 154.72 | ng/ul | 0.01 |
| 13) 4-Methylphenol-d8          | 8.46  | 113 | 2025707 | 138.43 | ng/ul | 0.01 |
| 19) Nitrobenzene-d5            | 8.90  | 128 | 976850  | 159.29 | ng/ul | 0.01 |
| 22) 2-Nitrophenol-d4           | 9.62  | 143 | 1028605 | 161.37 | ng/ul | 0.01 |
| 26) 2,4-Dichlorophenol-d3      | 10.16 | 165 | 2028544 | 158.08 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.67 | 131 | 1879557 | 132.71 | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.81 | 166 | 5398438 | 149.29 | ng/ul | 0.01 |
| 47) Acenaphthylene-d8          | 14.08 | 160 | 7008233 | 160.24 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.60 | 143 | 1029695 | 151.52 | ng/ul | 0.03 |
| 58) Fluorene-d10               | 15.39 | 176 | 4627441 | 147.74 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.51 | 200 | 837995  | 155.50 | ng/ul | 0.02 |
| 71) Anthracene-d10             | 17.24 | 188 | 6908677 | 151.48 | ng/ul | 0.01 |
| 79) Pyrene-d10                 | 19.53 | 212 | 7542893 | 163.06 | ng/ul | 0.01 |
| 90) Benzo(a)pyrene-d12         | 23.44 | 264 | 8666588 | 176.24 | ng/ul | 0.02 |

## Target Compounds

|                                |       |     |         |               | Qvalue |
|--------------------------------|-------|-----|---------|---------------|--------|
| 2) 1,4-Dioxane                 | 3.30  | 88  | 289324  | 60.826 ng/uL# | 74     |
| 4) Benzaldehyde                | 6.89  | 77  | 680323  | 63.518 ng/ul  | 93     |
| 6) Phenol                      | 6.95  | 94  | 2600984 | 139.790 ng/ul | 99     |
| 8) Bis(2-Chloroethyl)ether     | 7.19  | 93  | 1973097 | 134.699 ng/ul | 93     |
| 10) 2-Chlorophenol             | 7.32  | 128 | 2163954 | 157.030 ng/ul | 98     |
| 11) 2-Methylphenol             | 8.19  | 108 | 1962009 | 139.074 ng/ul | 97     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.28  | 45  | 3092628 | 130.256 ng/ul | 94     |
| 14) Acetophenone               | 8.57  | 105 | 3042709 | 136.034 ng/ul | 97     |
| 15) N-Nitroso-di-n-propylamine | 8.58  | 70  | 1547025 | 128.892 ng/ul | 92     |
| 16) 4-Methylphenol             | 8.53  | 108 | 2044132 | 132.424 ng/ul | 98     |
| 17) Hexachloroethane           | 8.82  | 117 | 859580  | 163.167 ng/ul | 88     |
| 20) Nitrobenzene               | 8.95  | 77  | 2250316 | 156.054 ng/ul | 90     |
| 21) Isophorone                 | 9.48  | 82  | 4227228 | 145.597 ng/ul | 97     |
| 23) 2-Nitrophenol              | 9.65  | 139 | 1112605 | 166.103 ng/ul | 93     |
| 24) 2,4-Dimethylphenol         | 9.72  | 107 | 2283311 | 159.313 ng/ul | 93     |
| 25) Bis(2-Chloroethoxy)methane | 9.96  | 93  | 2581453 | 136.360 ng/ul | 97     |
| 27) 2,4-Dichlorophenol         | 10.18 | 162 | 1922376 | 155.368 ng/ul | 97     |
| 28) Naphthalene                | 10.58 | 128 | 6401212 | 152.805 ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.69 | 127 | 1890069 | 140.459 ng/ul | 97     |
| 31) Hexachlorobutadiene        | 10.87 | 225 | 1261887 | 179.691 ng/ul | 97     |
| 32) Caprolactam                | 11.50 | 113 | 347430  | 89.201 ng/ul  | 88     |
| 33) 4-Chloro-3-methylphenol    | 11.83 | 107 | 2034508 | 155.092 ng/ul | 95     |
| 34) 2-Methylnaphthalene        | 12.20 | 142 | 4539949 | 145.935 ng/ul | 99     |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM091018\  
 Data File : BM016710.D  
 Acq On : 10 Sep 2018 14:52  
 Operator : SJ/JU  
 Sample : SSTD16052  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD16052

Quant Time: Sep 10 15:23:19 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 10 13:31:33 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|---------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.20 | 142  | 4539949  | 153.224 | ng/ul# | 94       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.57 | 216  | 2310511  | 182.099 | ng/ul# | 97       |
| 38) Hexachlorocyclopentadiene  | 12.55 | 237  | 1588276  | 223.327 | ng/ul  | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.81 | 196  | 1445472  | 167.326 | ng/ul  | 98       |
| 40) 2,4,5-Trichlorophenol      | 12.89 | 196  | 1564371  | 178.598 | ng/ul  | 98       |
| 41) 1,1'-Biphenyl              | 13.22 | 154  | 5505329  | 154.672 | ng/ul# | 98       |
| 42) 2-Chloronaphthalene        | 13.26 | 162  | 4312440  | 161.022 | ng/ul  | 96       |
| 43) 2-Nitroaniline             | 13.47 | 65   | 1268655  | 152.517 | ng/ul  | 97       |
| 45) Dimethylphthalate          | 13.86 | 163  | 5194797  | 150.047 | ng/ul  | 99       |
| 46) 2,6-Dinitrotoluene         | 13.97 | 165  | 1100411  | 179.969 | ng/ul  | 97       |
| 48) Acenaphthylene             | 14.11 | 152  | 6548299  | 148.768 | ng/ul  | 100      |
| 49) 3-Nitroaniline             | 14.30 | 138  | 943954   | 125.201 | ng/ul# | 96       |
| 50) Acenaphthene               | 14.45 | 153  | 4672849  | 157.117 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 14.50 | 184  | 537698   | 142.334 | ng/ul# | 90       |
| 53) 4-Nitrophenol              | 14.62 | 109  | 748822   | 171.447 | ng/ul# | 74       |
| 54) Dibenzofuran               | 14.79 | 168  | 6256999  | 150.404 | ng/ul  | 99       |
| 55) 2,4-Dinitrotoluene         | 14.76 | 165  | 1570097  | 174.682 | ng/ul  | 97       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.02 | 232  | 1341208  | 168.008 | ng/ul# | 86       |
| 57) Diethylphthalate           | 15.24 | 149  | 5327257  | 147.860 | ng/ul  | 98       |
| 59) Fluorene                   | 15.44 | 166  | 5161017  | 148.025 | ng/ul  | 100      |
| 60) 4-Chlorophenyl-phenylether | 15.44 | 204  | 2635601  | 161.335 | ng/ul  | 93       |
| 61) 4-Nitroaniline             | 15.47 | 138  | 1223450  | 151.488 | ng/ul  | 88       |
| 64) 4,6-Dinitro-2-methylphenol | 15.53 | 198  | 892013   | 161.533 | ng/ul  | 100      |
| 65) N-Nitrosodiphenylamine     | 15.66 | 169  | 4489198  | 159.127 | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.34 | 248  | 1689295  | 178.579 | ng/ul  | 97       |
| 67) Hexachlorobenzene          | 16.44 | 284  | 1784715  | 171.391 | ng/ul# | 94       |
| 68) Atrazine                   | 16.62 | 200  | 1639940  | 167.773 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 16.78 | 266  | 1123629  | 168.052 | ng/ul  | 97       |
| 70) Phenanthrene               | 17.18 | 178  | 7927283  | 153.558 | ng/ul  | 99       |
| 72) Anthracene                 | 17.27 | 178  | 8080013  | 152.532 | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.18 | 216  | 2419256  | 196.025 | ng/uL  | 98       |
| 74) Pentachlorobenzene         | 14.70 | 250  | 2187269  | 174.743 | ng/uL  | 99       |
| 75) Carbazole                  | 17.54 | 167  | 7311370  | 161.780 | ng/ul  | 97       |
| 76) Di-n-butylphthalate        | 18.13 | 149  | 8878334  | 149.568 | ng/ul  | 100      |
| 77) Fluoranthene               | 19.20 | 202  | 9268299  | 174.551 | ng/ul# | 89       |
| 80) Pyrene                     | 19.56 | 202  | 9149061  | 158.728 | ng/ul# | 84       |
| 81) Butylbenzylphthalate       | 20.48 | 149  | 4139602  | 156.168 | ng/ul  | 95       |
| 82) 3,3'-Dichlorobenzidine     | 21.25 | 252  | 2983270  | 166.280 | ng/ul# | 98       |
| 83) Benzo(a)anthracene         | 21.31 | 228  | 9261914  | 170.269 | ng/ul  | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 6057000  | 148.791 | ng/ul  | 99       |
| 85) Chrysene                   | 21.37 | 228  | 8585257  | 174.095 | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.16 | 149  | 10168640 | 138.861 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 22.91 | 252  | 9531512  | 153.551 | ng/ul# | 95       |
| 89) Benzo(k)fluoranthene       | 22.96 | 252  | 9319840  | 156.333 | ng/ul# | 95       |
| 91) Benzo(a)pyrene             | 23.49 | 252  | 9279121  | 158.372 | ng/ul# | 95       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.86 | 276  | 11279828 | 175.012 | ng/ul# | 86       |
| 93) Dibenzo(a,h)anthracene     | 25.87 | 278  | 9313250  | 170.414 | ng/ul# | 94       |
| 94) Benzo(g,h,i)perylene       | 26.55 | 276  | 9369874  | 179.514 | ng/ul# | 91       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM091018\  
Data File : BM016710.D  
Acq On : 10 Sep 2018 14:52  
Operator : SJ/JU  
Sample : SSTD16052  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD16052

Quant Time: Sep 10 15:23:19 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Sep 10 13:31:33 2018  
Response via : Initial Calibration

| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |

```
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091018\  
Data File : BM016710.D  
Acq On    : 10 Sep 2018  14:52  
Operator  : SJ/JU  
Sample    : SSTD16052  
Misc      :  
ALS Vial  : 7      Sample Multiplier: 1
```