

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM072715\  
 Data File : BM002062.D  
 Acq On : 27 Jul 2015 15:31  
 Operator : UM/HP  
 Sample : SSTD16077  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD16077

Manual Integrations  
 APPROVED

apatel  
 7/29/2015 10:42:49 AM

Quant Time: Jul 27 16:19:08 2015  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 27 14:35:16 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.60  | 152  | 68596    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.39 | 136  | 270252   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.26 | 164  | 156938   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.02 | 188  | 345667   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.22 | 240  | 403704   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.41 | 264  | 393133   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |        |       |      |
|--------------------------------|-------|-----|--------|--------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d     | 0.00   | ng/uL |      |
| 5) Phenol-d5                   | 6.79  | 99  | 821183 | 155.97 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 437684 | 159.03 | ng/ul | 0.01 |
| 9) 2-Chlorophenol-d4           | 0.00  | 132 | 0d     | 0.00   | ng/ul |      |
| 13) 4-Methylphenol-d8          | 8.33  | 113 | 659171 | 164.85 | 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         | 10.54 | 131 | 673151 | 157.66 | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 0.00  | 166 | 0d     | 0.00   | ng/ul |      |
| 47) Acenaphthylene-d8          | 0.00  | 160 | 0d     | 0.00   | ng/ul |      |
| 52) 4-Nitrophenol-d4           | 14.51 | 143 | 334743 | 183.46 | ng/ul | 0.04 |
| 58) Fluorene-d10               | 0.00  | 176 | 0d     | 0.00   | ng/ul |      |
| 63) 4,6-Dinitro-2-methylphenol | 15.42 | 200 | 380169 | 177.29 | ng/ul | 0.02 |
| 71) Anthracene-d10             | 0.00  | 188 | 0d     | 0.00   | ng/ul |      |
| 79) Pyrene-d10                 | 0.00  | 212 | 0d     | 0.00   | ng/ul |      |
| 90) Benzo(a)pyrene-d12         | 0.00  | 264 | 0d     | 0.00   | ng/ul |      |

## Target Compounds

|                                |       |     |         |        |        | Qvalue |
|--------------------------------|-------|-----|---------|--------|--------|--------|
| 4) Benzaldehyde                | 6.75  | 77  | 339781  | 121.47 | ng/ul  | 98     |
| 6) Phenol                      | 6.82  | 94  | 835109  | 157.82 | ng/ul  | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.04  | 93  | 611097  | 157.63 | ng/ul  | 96     |
| 11) 2-Methylphenol             | 8.06  | 108 | 639543  | 163.14 | ng/ul  | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.14  | 45  | 618616  | 176.03 | ng/ul  | 94     |
| 14) Acetophenone               | 8.44  | 105 | 1031080 | 165.66 | ng/ul  | 98     |
| 16) 4-Methylphenol             | 8.40  | 108 | 685269  | 164.03 | ng/ul  | 94     |
| 30) 4-Chloroaniline            | 10.56 | 127 | 703184  | 153.52 | ng/ul  | 99     |
| 32) Caprolactam                | 11.37 | 113 | 197218m | 159.70 | ng/ul  |        |
| 38) Hexachlorocyclopentadiene  | 12.41 | 237 | 568358  | 173.10 | ng/ul  | 98     |
| 49) 3-Nitroaniline             | 14.19 | 138 | 337238  | 171.11 | ng/ul  | 95     |
| 51) 2,4-Dinitrophenol          | 14.41 | 184 | 277074  | 203.89 | ng/ul  | 97     |
| 53) 4-Nitrophenol              | 14.52 | 109 | 279584  | 183.18 | ng/ul  | 98     |
| 61) 4-Nitroaniline             | 15.37 | 138 | 380232  | 185.35 | ng/ul  | 98     |
| 64) 4,6-Dinitro-2-methylphenol | 15.43 | 198 | 392815  | 175.28 | ng/ul  | 95     |
| 68) Atrazine                   | 16.50 | 200 | 613571  | 159.75 | ng/ul  | 100    |
| 69) Pentachlorophenol          | 16.67 | 266 | 383307  | 182.27 | ng/ul  | 97     |
| 75) Carbazole                  | 17.43 | 167 | 2535662 | 162.37 | ng/ul  | 99     |
| 77) Fluoranthene               | 19.08 | 202 | 3346593 | 158.85 | ng/ul# | 96     |
| 82) 3,3'-Dichlorobenzidine     | 21.14 | 252 | 1153624 | 166.66 | ng/ul  | 99     |
| 87) Di-n-octyl phthalate       | 22.00 | 149 | 3622015 | 165.24 | ng/ul  | 100    |

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM072715\  
Data File : BM002062.D  
Acq On : 27 Jul 2015 15:31  
Operator : UM/HP  
Sample : SSTD16077  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD16077

Manual Integrations  
APPROVED

apatel  
7/29/2015 10:42:49 AM

Quant Time: Jul 27 16:19:08 2015  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 27 14:35:16 2015  
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:\HPCHEM1\BNA\_M\DATA\BM072715\  
Data File : BM002062.D  
Acq On : 27 Jul 2015 15:31  
Operator : UM/HP  
Sample : SSTD16077  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD16077

Manual Integrations  
APPROVED

apatel  
7/29/2015 10:42:49 AM

Quant Time: Jul 27 16:19:08 2015  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 27 14:35:16 2015  
Response via : Initial Calibration

