

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030719\  
 Data File : BN004671.D  
 Acq On : 08 Mar 2019 11:29  
 Operator : JU/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD02025

Manual Integrations  
 APPROVED

Quant Time: Mar 08 14:29:38 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030719MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 08 00:36:35 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.72  | 152  | 36454    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.51 | 136  | 170686   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.36 | 164  | 101575   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.12 | 188  | 229192   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.32 | 240  | 194919   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.57 | 264  | 175587   | 20.00 | ng/ul | 0.00     |

Sohil

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 4651   | 7.20  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.89  | 99  | 57682  | 19.15 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.06  | 67  | 29473  | 18.76 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.25  | 132 | 51815  | 19.82 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.42  | 113 | 50671  | 19.34 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.88  | 128 | 25749  | 19.79 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.60  | 143 | 30175  | 20.42 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.13 | 165 | 56878  | 20.19 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.65 | 131 | 64320  | 19.04 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.77 | 166 | 164723 | 18.76 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.06 | 160 | 209097 | 19.78 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.57 | 143 | 29572  | 18.14 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.36 | 176 | 144244 | 19.32 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.49 | 200 | 26116  | 16.51 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.21 | 188 | 220821 | 19.78 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.52 | 212 | 227060 | 23.84 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.43 | 264 | 185452 | 19.67 | ng/ul | 0.00 |

3/8/2019 3:59:41 PM

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.26  | 88  | 5084   | 6.383  | ng/uL 99  |
| 4) Benzaldehyde                | 6.88  | 77  | 38386  | 19.681 | ng/ul 97  |
| 6) Phenol                      | 6.91  | 94  | 58729  | 19.157 | ng/ul 99  |
| 8) Bis(2-Chloroethyl)ether     | 7.15  | 93  | 42537  | 18.851 | ng/ul 98  |
| 10) 2-Chlorophenol             | 7.28  | 128 | 51420  | 19.497 | ng/ul 99  |
| 11) 2-Methylphenol             | 8.16  | 108 | 47453  | 19.274 | ng/ul 99  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.25  | 45  | 49141m | 18.882 | ng/ul     |
| 14) Acetophenone               | 8.55  | 105 | 76424  | 19.453 | ng/ul 99  |
| 15) N-Nitroso-di-n-propylamine | 8.53  | 70  | 35003  | 18.872 | ng/ul 98  |
| 16) 4-Methylphenol             | 8.49  | 108 | 52371  | 19.133 | ng/ul 97  |
| 17) Hexachloroethane           | 8.79  | 117 | 19629  | 19.872 | ng/ul 93  |
| 20) Nitrobenzene               | 8.93  | 77  | 52300  | 19.577 | ng/ul 99  |
| 21) Isophorone                 | 9.44  | 82  | 103820 | 18.757 | ng/ul 99  |
| 23) 2-Nitrophenol              | 9.63  | 139 | 31941  | 19.991 | ng/ul 99  |
| 24) 2,4-Dimethylphenol         | 9.68  | 107 | 60343  | 19.569 | ng/ul 99  |
| 25) Bis(2-Chloroethoxy)methane | 9.92  | 93  | 63298  | 18.787 | ng/ul 100 |
| 27) 2,4-Dichlorophenol         | 10.15 | 162 | 55393  | 19.984 | ng/ul 99  |
| 28) Naphthalene                | 10.56 | 128 | 173994 | 19.564 | ng/ul 99  |
| 30) 4-Chloroaniline            | 10.68 | 127 | 66134  | 19.258 | ng/ul 100 |
| 31) Hexachlorobutadiene        | 10.82 | 225 | 37519  | 20.366 | ng/ul 98  |
| 32) Caprolactam                | 11.45 | 113 | 17141  | 18.255 | ng/ul 97  |
| 33) 4-Chloro-3-methylphenol    | 11.79 | 107 | 57550  | 19.363 | ng/ul 99  |
| 34) 2-Methylnaphthalene        | 12.17 | 142 | 129199 | 19.388 | ng/ul 99  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030719\  
 Data File : BN004671.D  
 Acq On : 08 Mar 2019 11:29  
 Operator : JU/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD02025

Manual Integrations  
 APPROVED

Quant Time: Mar 08 14:29:38 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030719MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 08 00:36:35 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.39 | 142  | 121056   | 19.344 | ng/ul | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.54 | 216  | 74183    | 20.769 | ng/ul | 98       |
| 38) Hexachlorocyclopentadiene  | 12.51 | 237  | 34120    | 13.867 | ng/ul | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.79 | 196  | 47170    | 20.502 | ng/ul | 98       |
| 40) 2,4,5-Trichlorophenol      | 12.86 | 196  | 51125    | 20.299 | ng/ul | 96       |
| 41) 1,1'-Biphenyl              | 13.19 | 154  | 168015   | 19.661 | ng/ul | 99       |
| 42) 2-Chloronaphthalene        | 13.23 | 162  | 130921   | 20.150 | ng/ul | 99       |
| 43) 2-Nitroaniline             | 13.45 | 65   | 31263    | 19.668 | ng/ul | 99       |
| 45) Dimethylphthalate          | 13.82 | 163  | 162918   | 18.906 | ng/ul | 100      |
| 46) 2,6-Dinitrotoluene         | 13.95 | 165  | 34663    | 19.866 | ng/ul | 95       |
| 48) Acenaphthylene             | 14.09 | 152  | 203487   | 19.697 | ng/ul | 100      |
| 49) 3-Nitroaniline             | 14.28 | 138  | 33943    | 20.063 | ng/ul | 99       |
| 50) Acenaphthene               | 14.43 | 153  | 138104   | 19.556 | ng/ul | 100      |
| 51) 2,4-Dinitrophenol          | 14.49 | 184  | 16122    | 14.257 | ng/ul | 98       |
| 53) 4-Nitrophenol              | 14.59 | 109  | 22180    | 18.727 | ng/ul | 97       |
| 54) Dibenzofuran               | 14.76 | 168  | 197029   | 19.424 | ng/ul | 98       |
| 55) 2,4-Dinitrotoluene         | 14.74 | 165  | 50036    | 19.414 | ng/ul | 98       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.99 | 232  | 46196    | 19.818 | ng/ul | 99       |
| 57) Diethylphthalate           | 15.19 | 149  | 164931   | 18.887 | ng/ul | 100      |
| 59) Fluorene                   | 15.42 | 166  | 157812   | 19.390 | ng/ul | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.41 | 204  | 83771    | 19.387 | ng/ul | 100      |
| 61) 4-Nitroaniline             | 15.45 | 138  | 37557    | 18.912 | ng/ul | 100      |
| 64) 4,6-Dinitro-2-methylphenol | 15.50 | 198  | 27407    | 16.428 | ng/ul | 99       |
| 65) N-Nitrosodiphenylamine     | 15.63 | 169  | 138044   | 19.777 | ng/ul | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.30 | 248  | 56244    | 19.997 | ng/ul | 97       |
| 67) Hexachlorobenzene          | 16.41 | 284  | 64698    | 20.041 | ng/ul | 98       |
| 68) Atrazine                   | 16.57 | 200  | 52995    | 19.877 | ng/ul | 99       |
| 69) Pentachlorophenol          | 16.76 | 266  | 37549    | 18.663 | ng/ul | 98       |
| 70) Phenanthrene               | 17.16 | 178  | 248729   | 19.516 | ng/ul | 99       |
| 72) Anthracene                 | 17.25 | 178  | 255845   | 19.648 | ng/ul | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.15 | 216  | 72386    | 21.278 | ng/uL | 99       |
| 74) Pentachlorobenzene         | 14.67 | 250  | 74946    | 20.663 | ng/uL | 98       |
| 75) Carbazole                  | 17.52 | 167  | 224982   | 19.895 | ng/ul | 100      |
| 76) Di-n-butylphthalate        | 18.07 | 149  | 280602   | 20.009 | ng/ul | 100      |
| 77) Fluoranthene               | 19.18 | 202  | 284042   | 19.330 | ng/ul | 97       |
| 80) Pvrene                     | 19.55 | 202  | 281090   | 23.786 | ng/ul | 99       |
| 81) Butylbenzylphthalate       | 20.45 | 149  | 123886   | 25.534 | ng/ul | 99       |
| 82) 3,3'-Dichlorobenzidine     | 21.23 | 252  | 95043    | 20.654 | ng/ul | 100      |
| 83) Benzo(a)anthracene         | 21.30 | 228  | 241223   | 19.672 | ng/ul | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.22 | 149  | 174203   | 24.707 | ng/ul | 100      |
| 85) Chrysene                   | 21.35 | 228  | 219117   | 19.243 | ng/ul | 99       |
| 87) Di-n-octyl phthalate       | 22.10 | 149  | 262557   | 23.173 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.89 | 252  | 213123   | 19.556 | ng/ul | 99       |
| 89) Benzo(k)fluoranthene       | 22.94 | 252  | 200888   | 18.746 | ng/ul | 99       |
| 91) Benzo(a)pyrene             | 23.48 | 252  | 198570   | 19.368 | ng/ul | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.86 | 276  | 238776   | 21.381 | ng/ul | 98       |
| 93) Dibenzo(a,h)anthracene     | 25.87 | 278  | 200958   | 21.387 | ng/ul | 98       |
| 94) Benzo(a,h,i)perylene       | 26.56 | 276  | 200059   | 22.027 | ng/ul | 97       |

Sohil

3/8/2019 3:59:41 PM

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030719\  
Data File : BN004671.D  
Acq On : 08 Mar 2019 11:29  
Operator : JU/SJ  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD02025

Manual Integrations  
APPROVED

Quant Time: Mar 08 14:29:38 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030719MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Mar 08 00:36:35 2019  
Response via : Initial Calibration

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

Sohil

3/8/2019 3:59:41 PM

