

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031721\  
 Data File : BN013957.D  
 Acq On : 16 Mar 2021 16:29  
 Operator : CG/JU  
 Sample : SSTD01055  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010055

Manual Integrations  
 APPROVED

mohammad  
 3/17/2021 10:52:20 AM

Quant Time: Mar 16 17:43:28 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN031721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 16 17:38:29 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.71  | 152  | 137100   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.49 | 136  | 571342   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.35 | 164  | 337257   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.09 | 188  | 688550   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.27 | 240  | 683122   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.50 | 264  | 663995   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.21  | 96  | 14652  | 4.36  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.61  | 84  | 84734  | 9.65  | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.87  | 99  | 108765 | 10.09 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.05  | 67  | 76261  | 12.03 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.24  | 132 | 87723  | 9.30  | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.40  | 113 | 84204  | 9.84  | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.86  | 128 | 37772  | 8.42  | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.57  | 143 | 33471  | 7.27  | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.11 | 165 | 80281  | 9.11  | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.62 | 131 | 121182 | 9.74  | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.76 | 166 | 240483 | 9.61  | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.04 | 160 | 294311 | 9.25  | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.54 | 143 | 35964  | 7.67  | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.34 | 176 | 219618 | 10.10 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.46 | 200 | 24658  | 6.15  | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.19 | 188 | 329757 | 9.91  | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.48 | 212 | 363823 | 8.96  | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.35 | 264 | 346290 | 10.00 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.25  | 88  | 14738  | 4.236  | ng/uL 96  |
| 5) Pyridine                    | 3.63  | 79  | 91784  | 10.325 | ng/ul 94  |
| 6) Benzaldehyde                | 6.85  | 77  | 71761  | 16.310 | ng/ul 98  |
| 8) Phenol                      | 6.90  | 94  | 126462 | 11.358 | ng/ul 98  |
| 10) Bis(2-Chloroethyl)ether    | 7.14  | 93  | 101471 | 11.086 | ng/ul 99  |
| 12) 2-Chlorophenol             | 7.27  | 128 | 97968  | 10.139 | ng/ul 99  |
| 13) 2-Methylphenol             | 8.15  | 108 | 87765  | 10.463 | ng/ul 96  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.23  | 45  | 160224 | 12.537 | ng/ul 98  |
| 16) Acetophenone               | 8.52  | 105 | 152744 | 11.818 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.51  | 70  | 70539  | 11.995 | ng/ul 98  |
| 18) 4-Methylphenol             | 8.47  | 108 | 99810  | 10.989 | ng/ul 97  |
| 19) Hexachloroethane           | 8.78  | 117 | 40668  | 10.105 | ng/ul 99  |
| 22) Nitrobenzene               | 8.90  | 77  | 109378 | 10.750 | ng/ul 100 |
| 23) Isophorone                 | 9.42  | 82  | 177272 | 10.139 | ng/ul 99  |
| 25) 2-Nitrophenol              | 9.61  | 139 | 41004  | 7.844  | ng/ul 99  |
| 26) 2,4-Dimethylphenol         | 9.67  | 107 | 106466 | 10.619 | ng/ul 99  |
| 27) Bis(2-Chloroethoxy)methane | 9.91  | 93  | 130403 | 10.697 | ng/ul 99  |
| 29) 2,4-Dichlorophenol         | 10.13 | 162 | 88072  | 9.927  | ng/ul 97  |
| 30) Naphthalene                | 10.54 | 128 | 335575 | 10.404 | ng/ul 100 |
| 32) 4-Chloroaniline            | 10.65 | 127 | 130054 | 10.822 | ng/ul 97  |
| 33) Hexachlorobutadiene        | 10.83 | 225 | 58107  | 10.033 | ng/ul 93  |
| 34) Caprolactam                | 11.39 | 113 | 17288m | 7.040  | ng/ul     |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031721\  
 Data File : BN013957.D  
 Acq On : 16 Mar 2021 16:29  
 Operator : CG/JU  
 Sample : SSTD01055  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010055

Manual Integrations  
 APPROVED

mohammad  
 3/17/2021 10:52:20 AM

Quant Time: Mar 16 17:43:28 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN031721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 16 17:38:29 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 4-Chloro-3-methylphenol    | 11.77 | 107  | 84999    | 9.773  | ng/ul | 99       |
| 36) 2-Methylnaphthalene        | 12.16 | 142  | 223793   | 10.378 | ng/ul | 97       |
| 37) 1-Methylnaphthalene        | 12.37 | 142  | 225683   | 10.930 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.53 | 216  | 112944   | 10.424 | ng/ul | 97       |
| 40) Hexachlorocyclopentadiene  | 12.51 | 237  | 59416    | 8.943  | ng/ul | 98       |
| 41) 2,4,6-Trichlorophenol      | 12.77 | 196  | 56558    | 8.656  | ng/ul | 96       |
| 42) 2,4,5-Trichlorophenol      | 12.83 | 196  | 65065    | 9.109  | ng/ul | 96       |
| 43) 1,1'-Biphenyl              | 13.17 | 154  | 288367   | 10.135 | ng/ul | 97       |
| 44) 2-Chloronaphthalene        | 13.22 | 162  | 225406   | 10.022 | ng/ul | 100      |
| 45) 2-Nitroaniline             | 13.42 | 65   | 45081    | 8.773  | ng/ul | 96       |
| 47) Dimethylphthalate          | 13.80 | 163  | 257083   | 9.878  | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene         | 13.92 | 165  | 39328    | 7.723  | ng/ul | 94       |
| 50) Acenaphthylene             | 14.07 | 152  | 325726   | 9.934  | ng/ul | 96       |
| 51) 3-Nitroaniline             | 14.25 | 138  | 44050    | 9.431  | ng/ul | 99       |
| 52) Acenaphthene               | 14.41 | 153  | 238933   | 10.212 | ng/ul | 97       |
| 53) 2,4-Dinitrophenol          | 14.45 | 184  | 15391    | 5.924  | ng/ul | 97       |
| 55) 4-Nitrophenol              | 14.55 | 109  | 32489    | 9.633  | ng/ul | 97       |
| 56) Dibenzofuran               | 14.75 | 168  | 330481   | 10.547 | ng/ul | 94       |
| 57) 2,4-Dinitrotoluene         | 14.71 | 165  | 58394    | 8.207  | ng/ul | 93       |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.97 | 232  | 48002    | 8.674  | ng/ul | 95       |
| 59) Diethylphthalate           | 15.17 | 149  | 245853   | 9.780  | ng/ul | 99       |
| 61) Fluorene                   | 15.40 | 166  | 269487   | 10.901 | ng/ul | 98       |
| 62) 4-Chlorophenyl-phenylether | 15.39 | 204  | 131187   | 10.815 | ng/ul | 93       |
| 63) 4-Nitroaniline             | 15.41 | 138  | 46386    | 11.049 | ng/ul | 99       |
| 66) 4,6-Dinitro-2-methylphenol | 15.47 | 198  | 27604    | 6.346  | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine     | 15.61 | 169  | 221307   | 9.703  | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether  | 16.29 | 248  | 72744    | 9.369  | ng/ul | 97       |
| 69) Hexachlorobenzene          | 16.40 | 284  | 86437    | 9.722  | ng/ul | 97       |
| 70) Atrazine                   | 16.56 | 200  | 61075    | 8.183  | ng/ul | 99       |
| 71) Pentachlorophenol          | 16.74 | 266  | 32289    | 6.805  | ng/ul | 92       |
| 72) Phenanthrene               | 17.13 | 178  | 414662   | 10.086 | ng/ul | 97       |
| 74) Anthracene                 | 17.22 | 178  | 426615   | 10.337 | ng/ul | 98       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.14 | 216  | 114064   | 9.757  | ng/uL | 95       |
| 76) Pentachlorobenzene         | 14.66 | 250  | 112458   | 10.112 | ng/uL | 97       |
| 77) Carbazole                  | 17.49 | 167  | 358071   | 10.383 | ng/ul | 99       |
| 78) Di-n-butylphthalate        | 18.07 | 149  | 341746   | 8.717  | ng/ul | 100      |
| 80) Fluoranthene               | 19.15 | 202  | 445430   | 8.447  | ng/ul | 99       |
| 82) Pyrene                     | 19.51 | 202  | 488856   | 9.030  | ng/ul | 97       |
| 83) Butylbenzylphthalate       | 20.42 | 149  | 118295   | 6.566  | ng/ul | 97       |
| 84) 3,3'-Dichlorobenzidine     | 21.19 | 252  | 114548   | 9.456  | ng/ul | 99       |
| 85) Benzo(a)anthracene         | 21.25 | 228  | 454339   | 10.012 | ng/ul | 97       |
| 86) Bis(2-ethylhexyl)phthalate | 21.19 | 149  | 198747   | 7.843  | ng/ul | 99       |
| 87) Chrysene                   | 21.30 | 228  | 436901   | 9.781  | ng/ul | 97       |
| 89) Di-n-octyl phthalate       | 22.07 | 149  | 308040   | 7.535  | ng/ul | 100      |
| 90) Benzo(b)fluoranthene       | 22.83 | 252  | 450631   | 10.370 | ng/ul | 96       |
| 91) Benzo(k)fluoranthene       | 22.87 | 252  | 448542   | 10.427 | ng/ul | 98       |
| 93) Benzo(a)pyrene             | 23.40 | 252  | 396287   | 9.925  | ng/ul | 96       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.73 | 276  | 496634   | 10.030 | ng/ul | 97       |
| 95) Dibenzo(a,h)anthracene     | 25.74 | 278  | 423739   | 10.244 | ng/ul | 98       |
| 96) Benzo(g,h,i)perylene       | 26.42 | 276  | 412107   | 10.013 | ng/ul | 99       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031721\  
Data File : BN013957.D  
Acq On : 16 Mar 2021 16:29  
Operator : CG/JU  
Sample : SSTD01055  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010055

Manual Integrations  
APPROVED

mohammad  
3/17/2021 10:52:20 AM

Quant Time: Mar 16 17:43:28 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN031721.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Mar 16 17:38:29 2021  
Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

|       |  |  |  |  |  |  |
|-------|--|--|--|--|--|--|
| ----- |  |  |  |  |  |  |
| ----- |  |  |  |  |  |  |

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

