

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030421\  
 Data File : BN013828.D  
 Acq On : 04 Mar 2021 12:17  
 Operator : CG/JU  
 Sample : SSTD04054  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD04054

Manual Integrations  
 APPROVED

mohammad  
 3/5/2021 2:44:49 PM

Quant Time: Mar 04 13:36:01 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030421MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 04 12:54:17 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.72  | 152  | 161223   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.51 | 136  | 646318   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.36 | 164  | 371919   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.10 | 188  | 751913   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.28 | 240  | 797443   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.51 | 264  | 799373   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 64411   | 17.16 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.89  | 99  | 516965  | 43.18 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.06  | 67  | 314513  | 45.64 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.26  | 132 | 402813  | 40.26 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.43  | 113 | 389894  | 40.90 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.87  | 128 | 182441  | 41.70 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.59  | 143 | 180504  | 39.99 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.13 | 165 | 371642  | 41.61 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.64 | 131 | 492694  | 45.91 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.77 | 166 | 1006223 | 40.39 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.05 | 160 | 1366713 | 41.74 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.56 | 143 | 192255  | 39.78 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.36 | 176 | 877581  | 40.20 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.47 | 200 | 138792  | 34.51 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.20 | 188 | 1333430 | 40.90 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.49 | 212 | 1480603 | 39.08 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.37 | 264 | 1615242 | 41.38 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue    |
|--------------------------------|-------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.26  | 88  | 66704   | 17.571 | ng/uL 98  |
| 4) Benzaldehyde                | 6.87  | 77  | 308677  | 50.697 | ng/ul 97  |
| 6) Phenol                      | 6.92  | 94  | 547889  | 43.105 | ng/ul 98  |
| 8) Bis(2-Chloroethyl)ether     | 7.15  | 93  | 431310  | 43.467 | ng/ul 98  |
| 10) 2-Chlorophenol             | 7.29  | 128 | 428104  | 40.373 | ng/ul 99  |
| 11) 2-Methylphenol             | 8.16  | 108 | 398718  | 42.152 | ng/ul 97  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.26  | 45  | 648659  | 45.501 | ng/ul 98  |
| 14) Acetophenone               | 8.54  | 105 | 642838  | 42.639 | ng/ul 99  |
| 15) N-Nitroso-di-n-propylamine | 8.53  | 70  | 323919  | 46.376 | ng/ul 99  |
| 16) 4-Methylphenol             | 8.49  | 108 | 431032  | 41.604 | ng/ul 99  |
| 17) Hexachloroethane           | 8.80  | 117 | 180305  | 43.239 | ng/ul 97  |
| 20) Nitrobenzene               | 8.92  | 77  | 480985  | 47.125 | ng/ul 99  |
| 21) Isophorone                 | 9.44  | 82  | 856873  | 47.848 | ng/ul 99  |
| 23) 2-Nitrophenol              | 9.62  | 139 | 209228  | 40.245 | ng/ul 98  |
| 24) 2,4-Dimethylphenol         | 9.69  | 107 | 463621  | 43.503 | ng/ul 98  |
| 25) Bis(2-Chloroethoxy)methane | 9.93  | 93  | 544151  | 43.136 | ng/ul 100 |
| 27) 2,4-Dichlorophenol         | 10.15 | 162 | 373305  | 40.768 | ng/ul 96  |
| 28) Naphthalene                | 10.56 | 128 | 1314170 | 39.631 | ng/ul 100 |
| 30) 4-Chloroaniline            | 10.66 | 127 | 505054  | 45.912 | ng/ul 98  |
| 31) Hexachlorobutadiene        | 10.85 | 225 | 231875  | 41.243 | ng/ul 98  |
| 32) Caprolactam                | 11.42 | 113 | 112893m | 41.285 | ng/ul     |
| 33) 4-Chloro-3-methylphenol    | 11.79 | 107 | 391824  | 41.963 | ng/ul 98  |
| 34) 2-Methylnaphthalene        | 12.17 | 142 | 910101  | 39.283 | ng/ul 99  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030421\  
 Data File : BN013828.D  
 Acq On : 04 Mar 2021 12:17  
 Operator : CG/JU  
 Sample : SSTD04054  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 SSTD04054

Manual Integrations  
 APPROVED

mohammad  
 3/5/2021 2:44:49 PM

Quant Time: Mar 04 13:36:01 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030421MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 04 12:54:17 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.39 | 142  | 867696   | 38.782 | ng/ul | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.54 | 216  | 442007   | 42.286 | ng/ul | 99       |
| 38) Hexachlorocyclopentadiene  | 12.53 | 237  | 266358   | 41.609 | ng/ul | 98       |
| 39) 2,4,6-Trichlorophenol      | 12.78 | 196  | 273952   | 43.255 | ng/ul | 98       |
| 40) 2,4,5-Trichlorophenol      | 12.85 | 196  | 299537   | 42.059 | ng/ul | 97       |
| 41) 1,1'-Biphenyl              | 13.19 | 154  | 1101599  | 39.141 | ng/ul | 100      |
| 42) 2-Chloronaphthalene        | 13.23 | 162  | 860820   | 39.512 | ng/ul | 98       |
| 43) 2-Nitroaniline             | 13.43 | 65   | 251866   | 49.051 | ng/ul | 98       |
| 45) Dimethylphthalate          | 13.82 | 163  | 1038080  | 40.086 | ng/ul | 100      |
| 46) 2,6-Dinitrotoluene         | 13.94 | 165  | 205532   | 41.737 | ng/ul | 94       |
| 48) Acenaphthylene             | 14.08 | 152  | 1294706  | 40.395 | ng/ul | 99       |
| 49) 3-Nitroaniline             | 14.26 | 138  | 223038   | 44.977 | ng/ul | 100      |
| 50) Acenaphthene               | 14.42 | 153  | 922370   | 39.773 | ng/ul | 98       |
| 51) 2,4-Dinitrophenol          | 14.46 | 184  | 97990    | 34.966 | ng/ul | 97       |
| 53) 4-Nitrophenol              | 14.57 | 109  | 152965   | 43.108 | ng/ul | 94       |
| 54) Dibenzofuran               | 14.76 | 168  | 1253470  | 38.908 | ng/ul | 99       |
| 55) 2,4-Dinitrotoluene         | 14.72 | 165  | 289673   | 39.999 | ng/ul | 98       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.99 | 232  | 236588   | 42.294 | ng/ul | 98       |
| 57) Diethylphthalate           | 15.19 | 149  | 1036690  | 40.721 | ng/ul | 100      |
| 59) Fluorene                   | 15.41 | 166  | 1047859  | 40.618 | ng/ul | 97       |
| 60) 4-Chlorophenyl-phenylether | 15.41 | 204  | 497260   | 40.486 | ng/ul | 95       |
| 61) 4-Nitroaniline             | 15.43 | 138  | 236606   | 42.525 | ng/ul | 95       |
| 64) 4,6-Dinitro-2-methylphenol | 15.49 | 198  | 149156   | 35.877 | ng/ul | 95       |
| 65) N-Nitrosodiphenylamine     | 15.62 | 169  | 891501   | 41.058 | ng/ul | 98       |
| 66) 4-Bromophenyl-phenylether  | 16.30 | 248  | 290489   | 40.155 | ng/ul | 97       |
| 67) Hexachlorobenzene          | 16.41 | 284  | 340586   | 41.075 | ng/ul | 99       |
| 68) Atrazine                   | 16.57 | 200  | 305403   | 41.580 | ng/ul | 98       |
| 69) Pentachlorophenol          | 16.75 | 266  | 188577   | 40.184 | ng/ul | 97       |
| 70) Phenanthrene               | 17.15 | 178  | 1624344  | 40.248 | ng/ul | 99       |
| 72) Anthracene                 | 17.24 | 178  | 1668649  | 40.861 | ng/ul | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.15 | 216  | 435676   | 41.689 | ng/uL | 93       |
| 74) Pentachlorobenzene         | 14.68 | 250  | 414087   | 40.833 | ng/uL | 97       |
| 75) Carbazole                  | 17.50 | 167  | 1519710  | 42.731 | ng/ul | 99       |
| 76) Di-n-butylphthalate        | 18.08 | 149  | 1765942  | 44.252 | ng/ul | 99       |
| 77) Fluoranthene               | 19.16 | 202  | 1865984  | 43.575 | ng/ul | 96       |
| 80) Pvrene                     | 19.52 | 202  | 1951100  | 38.862 | ng/ul | 99       |
| 81) Butylbenzylphthalate       | 20.43 | 149  | 763512   | 42.814 | ng/ul | 99       |
| 82) 3,3'-Dichlorobenzidine     | 21.20 | 252  | 642352   | 45.120 | ng/ul | 96       |
| 83) Benzo(a)anthracene         | 21.27 | 228  | 1919774  | 40.067 | ng/ul | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.21 | 149  | 1242630  | 44.546 | ng/ul | 99       |
| 85) Chrysene                   | 21.32 | 228  | 1860855  | 39.195 | ng/ul | 97       |
| 87) Di-n-octyl phthalate       | 22.09 | 149  | 2042644  | 43.612 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.84 | 252  | 1959537  | 40.599 | ng/ul | 96       |
| 89) Benzo(k)fluoranthene       | 22.89 | 252  | 1928416  | 41.092 | ng/ul | 100      |
| 91) Benzo(a)pyrene             | 23.42 | 252  | 1775880  | 41.928 | ng/ul | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.76 | 276  | 2263020  | 41.676 | ng/ul | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.77 | 278  | 1891856  | 41.371 | ng/ul | 99       |
| 94) Benzo(a,h,i)perylene       | 26.45 | 276  | 1885713  | 41.267 | ng/ul | 97       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN030421\  
Data File : BN013828.D  
Acq On : 04 Mar 2021 12:17  
Operator : CG/JU  
Sample : SSTD04054  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD04054

Manual Integrations  
APPROVED

mohammad  
3/5/2021 2:44:49 PM

Quant Time: Mar 04 13:36:01 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030421MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Mar 04 12:54:17 2021  
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 N\DATA\BN030421\  
Data File : BN013828.D  
Acq On : 04 Mar 2021 12:17  
Operator : CG/JU  
Sample : SST04054  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SST04054

Manual Integrations  
APPROVED

mohammad  
3/5/2021 2:44:49 PM

Quant Time: Mar 04 13:36:01 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030421MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Mar 04 12:54:17 2021  
Response via : Initial Calibration

