

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
 Data File : BN014285.D  
 Acq On : 15 Apr 2021 15:07  
 Operator : CG/JU  
 Sample : SSTDICV020  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SICV057

Manual Integrations  
 APPROVED

mohammad  
 4/16/2021 11:30:54 AM

Quant Time: Apr 15 16:15:10 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 15 15:18:04 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.69  | 152  | 188779   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.47 | 136  | 766805   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.33 | 164  | 450808   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.08 | 188  | 940701   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.28 | 240  | 1011528  | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.51 | 264  | 1025862  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 34624  | 6.98  | ng/uL | 0.00  |
| 4) Pyridine-d5                 | 3.62  | 84  | 228788 | 17.41 | ng/ul | 0.00  |
| 7) Phenol-d5                   | 6.86  | 99  | 277330 | 17.57 | ng/ul | -0.01 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.03  | 67  | 165478 | 17.51 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4          | 7.23  | 132 | 230340 | 18.56 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8          | 8.39  | 113 | 213475 | 17.84 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5            | 8.84  | 128 | 105192 | 18.98 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4           | 9.56  | 143 | 105588 | 19.46 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3      | 10.09 | 165 | 217907 | 18.64 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4         | 10.60 | 131 | 299267 | 17.26 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6       | 13.75 | 166 | 602997 | 18.46 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8          | 14.02 | 160 | 712083 | 18.12 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4           | 14.53 | 143 | 112197 | 17.71 | ng/ul | 0.00  |
| 60) Fluorene-d10               | 15.33 | 176 | 517052 | 17.91 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylphenol | 15.44 | 200 | 83842  | 16.65 | ng/ul | 0.00  |
| 73) Anthracene-d10             | 17.18 | 188 | 790657 | 18.01 | ng/ul | 0.00  |
| 81) Pyrene-d10                 | 19.48 | 212 | 896310 | 18.25 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12         | 23.37 | 264 | 967016 | 17.77 | ng/ul | 0.00  |

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.25  | 88  | 36187  | 7.162  | ng/uL 98  |
| 5) Pyridine                    | 3.64  | 79  | 235290 | 17.423 | ng/ul 95  |
| 6) Benzaldehyde                | 6.84  | 77  | 179963 | 22.585 | ng/ul 99  |
| 8) Phenol                      | 6.89  | 94  | 300651 | 17.722 | ng/ul 99  |
| 10) Bis(2-Chloroethyl)ether    | 7.12  | 93  | 235911 | 17.682 | ng/ul 99  |
| 12) 2-Chlorophenol             | 7.26  | 128 | 244661 | 18.434 | ng/ul 98  |
| 13) 2-Methylphenol             | 8.13  | 108 | 218475 | 17.760 | ng/ul 97  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.23  | 45  | 263084 | 17.838 | ng/ul 99  |
| 16) Acetophenone               | 8.50  | 105 | 366307 | 18.217 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.50  | 70  | 176298 | 18.857 | ng/ul 98  |
| 18) 4-Methylphenol             | 8.46  | 108 | 239532 | 17.980 | ng/ul 100 |
| 19) Hexachloroethane           | 8.76  | 117 | 102076 | 18.297 | ng/ul 99  |
| 22) Nitrobenzene               | 8.88  | 77  | 269222 | 18.367 | ng/ul 100 |
| 23) Isophorone                 | 9.40  | 82  | 464170 | 18.829 | ng/ul 100 |
| 25) 2-Nitrophenol              | 9.59  | 139 | 124756 | 19.458 | ng/ul 99  |
| 26) 2,4-Dimethylphenol         | 9.65  | 107 | 259311 | 18.167 | ng/ul 98  |
| 27) Bis(2-Chloroethoxy)methane | 9.89  | 93  | 304287 | 18.118 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 10.12 | 162 | 224915 | 18.749 | ng/ul 98  |
| 30) Naphthalene                | 10.52 | 128 | 766247 | 18.029 | ng/ul 100 |
| 32) 4-Chloroaniline            | 10.63 | 127 | 312605 | 17.519 | ng/ul 99  |
| 33) Hexachlorobutadiene        | 10.81 | 225 | 147442 | 18.825 | ng/ul 92  |
| 34) Caprolactam                | 11.38 | 113 | 60761  | 18.172 | ng/ul 96  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
 Data File : BN014285.D  
 Acq On : 15 Apr 2021 15:07  
 Operator : CG/JU  
 Sample : SSTDICV020  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SICV057

Manual Integrations  
 APPROVED

mohammad  
 4/16/2021 11:30:54 AM

Quant Time: Apr 15 16:15:10 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 15 15:18:04 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 4-Chloro-3-methylphenol    | 11.76 | 107  | 223305   | 18.833 | ng/ul  | 99       |
| 36) 2-Methylnaphthalene        | 12.14 | 142  | 524154   | 18.101 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene        | 12.36 | 142  | 519467   | 18.014 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.51 | 216  | 271034   | 17.763 | ng/ul  | 96       |
| 40) Hexachlorocyclopentadiene  | 12.50 | 237  | 163231   | 17.227 | ng/ul  | 100      |
| 41) 2,4,6-Trichlorophenol      | 12.75 | 196  | 163654   | 18.557 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol      | 12.82 | 196  | 175173   | 17.899 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl              | 13.16 | 154  | 666153   | 17.967 | ng/ul  | 98       |
| 44) 2-Chloronaphthalene        | 13.20 | 162  | 520042   | 17.943 | ng/ul  | 100      |
| 45) 2-Nitroaniline             | 13.40 | 65   | 128088   | 19.063 | ng/ul  | 96       |
| 47) Dimethylphthalate          | 13.79 | 163  | 624529   | 18.421 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene         | 13.90 | 165  | 114311   | 19.226 | ng/ul  | 100      |
| 50) Acenaphthylene             | 14.05 | 152  | 768393   | 18.382 | ng/ul  | 100      |
| 51) 3-Nitroaniline             | 14.23 | 138  | 132259   | 18.053 | ng/ul  | 96       |
| 52) Acenaphthene               | 14.40 | 153  | 549343   | 18.016 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol          | 14.44 | 184  | 55272    | 15.926 | ng/ul  | 94       |
| 55) 4-Nitrophenol              | 14.54 | 109  | 94774    | 17.462 | ng/ul  | 99       |
| 56) Dibenzofuran               | 14.73 | 168  | 746458   | 17.591 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene         | 14.69 | 165  | 168194   | 19.582 | ng/ul  | 96       |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.96 | 232  | 145374   | 18.881 | ng/ul  | 99       |
| 59) Diethylphthalate           | 15.16 | 149  | 632931   | 18.250 | ng/ul  | 100      |
| 61) Fluorene                   | 15.38 | 166  | 626798   | 18.222 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenylether | 15.38 | 204  | 309084   | 18.247 | ng/ul  | 98       |
| 63) 4-Nitroaniline             | 15.40 | 138  | 135813   | 19.045 | ng/ul  | 100      |
| 66) 4,6-Dinitro-2-methylphenol | 15.46 | 198  | 89054    | 16.705 | ng/ul# | 99       |
| 67) N-Nitrosodiphenylamine     | 15.59 | 169  | 530538   | 18.126 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether  | 16.28 | 248  | 177111   | 17.797 | ng/ul  | 92       |
| 69) Hexachlorobenzene          | 16.39 | 284  | 212909   | 17.923 | ng/ul  | 96       |
| 70) Atrazine                   | 16.55 | 200  | 182039   | 17.501 | ng/ul  | 98       |
| 71) Pentachlorophenol          | 16.73 | 266  | 115660   | 16.745 | ng/ul  | 91       |
| 72) Phenanthrene               | 17.12 | 178  | 978373   | 17.814 | ng/ul  | 99       |
| 74) Anthracene                 | 17.22 | 178  | 993926   | 17.816 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.12 | 216  | 269527   | 17.399 | ng/uL  | 96       |
| 76) Pentachlorobenzene         | 14.65 | 250  | 274590   | 17.835 | ng/uL  | 93       |
| 77) Carbazole                  | 17.48 | 167  | 876513   | 17.368 | ng/ul  | 100      |
| 78) Di-n-butylphthalate        | 18.06 | 149  | 1012324  | 18.539 | ng/ul  | 99       |
| 80) Fluoranthene               | 19.14 | 202  | 1105192  | 17.812 | ng/ul  | 99       |
| 82) Pyrene                     | 19.51 | 202  | 1170270  | 17.713 | ng/ul  | 98       |
| 83) Butylbenzylphthalate       | 20.42 | 149  | 423544   | 19.524 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine     | 21.20 | 252  | 374771   | 20.716 | ng/ul  | 91       |
| 85) Benzo(a)anthracene         | 21.26 | 228  | 1176186  | 17.996 | ng/ul  | 96       |
| 86) Bis(2-ethylhexyl)phthalate | 21.21 | 149  | 700887   | 19.823 | ng/ul# | 98       |
| 87) Chrysene                   | 21.32 | 228  | 1143981  | 18.057 | ng/ul  | 95       |
| 89) Di-n-octyl phthalate       | 22.09 | 149  | 1131087  | 17.659 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 22.84 | 252  | 1211543  | 17.653 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene       | 22.89 | 252  | 1199840  | 17.562 | ng/ul  | 99       |
| 93) Benzo(a)pyrene             | 23.42 | 252  | 1095821  | 18.047 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.77 | 276  | 1432538m | 17.800 | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene     | 25.77 | 278  | 1219208  | 18.351 | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene       | 26.44 | 276  | 1146859  | 17.634 | ng/ul  | 100      |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
Data File : BN014285.D  
Acq On : 15 Apr 2021 15:07  
Operator : CG/JU  
Sample : SSTDICV020  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SICV057

Manual Integrations  
APPROVED

mohammad  
4/16/2021 11:30:54 AM

Quant Time: Apr 15 16:15:10 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Apr 15 15:18:04 2021  
Response via : Initial Calibration

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

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

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

