

Data File : BG046178.D  
Acq On : 6 Aug 2020 12:00  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD02040

Manual Integrations  
APPROVED

mohammad  
8/7/2020 1:49:53 PM

Quant Time: Aug 06 12:42:56 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG072320MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 06 10:37:54 2020  
Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.95  | 152  | 106237   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.75 | 136  | 431529   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.58 | 164  | 289803   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.32 | 188  | 657346   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.57 | 240  | 628143   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.67 | 264  | 678733   | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.42  | 96  | 16922  | 7.87  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.11  | 99  | 169012 | 20.27 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.28  | 67  | 96618  | 18.71 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.49  | 132 | 139735 | 20.94 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.66  | 113 | 137779 | 19.59 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.10  | 128 | 67131  | 22.03 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.83  | 143 | 76098  | 24.44 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.37 | 165 | 155205 | 22.02 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.88 | 131 | 173000 | 21.79 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.99 | 166 | 448463 | 19.84 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.27 | 160 | 546313 | 20.37 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.77 | 143 | 76106  | 20.62 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.57 | 176 | 414154 | 20.13 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.68 | 200 | 62498  | 19.00 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.42 | 188 | 621593 | 20.50 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.70 | 212 | 686675 | 20.67 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.46 | 264 | 749122 | 19.91 | ng/ul | 0.00 |

#### Target Compounds

|                                |       |     |        |        | Qvalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.45  | 88  | 20566  | 7.808  | ng/uL 96  |
| 4) Benzaldehyde                | 7.09  | 77  | 111287 | 18.724 | ng/ul 99  |
| 6) Phenol                      | 7.14  | 94  | 171133 | 19.929 | ng/ul 95  |
| 8) Bis(2-Chloroethyl)ether     | 7.37  | 93  | 137148 | 19.514 | ng/ul 96  |
| 10) 2-Chlorophenol             | 7.52  | 128 | 137694 | 20.355 | ng/ul 97  |
| 11) 2-Methylphenol             | 8.39  | 108 | 139638 | 20.181 | ng/ul 96  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.49  | 45  | 209491 | 17.636 | ng/ul# 94 |
| 14) Acetophenone               | 8.77  | 105 | 204531 | 19.709 | ng/ul 95  |
| 15) N-Nitroso-di-n-propylamine | 8.76  | 70  | 106056 | 18.888 | ng/ul 95  |
| 16) 4-Methylphenol             | 8.72  | 108 | 147763 | 19.932 | ng/ul 96  |
| 17) Hexachloroethane           | 9.04  | 117 | 56650  | 19.735 | ng/ul 92  |
| 20) Nitrobenzene               | 9.15  | 77  | 155375 | 19.849 | ng/ul 97  |
| 21) Isophorone                 | 9.67  | 82  | 304687 | 19.918 | ng/ul 98  |
| 23) 2-Nitrophenol              | 9.86  | 139 | 82125  | 23.558 | ng/ul 98  |
| 24) 2,4-Dimethylphenol         | 9.92  | 107 | 156615 | 20.274 | ng/ul 99  |
| 25) Bis(2-Chloroethoxy)methane | 10.16 | 93  | 188555 | 19.486 | ng/ul 99  |
| 27) 2,4-Dichlorophenol         | 10.40 | 162 | 149143 | 21.471 | ng/ul 97  |
| 28) Naphthalene                | 10.80 | 128 | 456882 | 19.783 | ng/ul 99  |
| 30) 4-Chloroaniline            | 10.90 | 127 | 169066 | 22.146 | ng/ul 89  |
| 31) Hexachlorobutadiene        | 11.10 | 225 | 114248 | 21.373 | ng/ul 96  |
| 32) Caprolactam                | 11.65 | 113 | 48756  | 20.821 | ng/ul 92  |
| 33) 4-Chloro-3-methylphenol    | 12.03 | 107 | 147368 | 20.691 | ng/ul 99  |
| 34) 2-Methylnaphthalene        | 12.41 | 142 | 347667 | 20.024 | ng/ul 97  |

Data File : BG046178.D  
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Sample : SSTDC0020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleID :  
SSTD02040

Manual Integrations  
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Quant Time: Aug 06 12:42:56 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG072320MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 06 10:37:54 2020  
Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.62 | 142  | 338014   | 19.712 | ng/uL | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.78 | 216  | 213156   | 21.275 | ng/uL | 96       |
| 38) Hexachlorocyclopentadiene  | 12.76 | 237  | 109016   | 19.004 | ng/uL | 95       |
| 39) 2,4,6-Trichlorophenol      | 13.01 | 196  | 135216   | 23.088 | ng/uL | 97       |
| 40) 2,4,5-Trichlorophenol      | 13.08 | 196  | 144107   | 22.635 | ng/uL | 97       |
| 41) 1,1'-Biphenyl              | 13.42 | 154  | 450390   | 20.087 | ng/uL | 97       |
| 42) 2-Chloronaphthalene        | 13.45 | 162  | 364149   | 20.355 | ng/uL | 98       |
| 43) 2-Nitroaniline             | 13.65 | 65   | 86232    | 20.448 | ng/uL | 96       |
| 45) Dimethylphthalate          | 14.03 | 163  | 446125   | 19.681 | ng/uL | 97       |
| 46) 2,6-Dinitrotoluene         | 14.15 | 165  | 97444    | 22.542 | ng/uL | 93       |
| 48) Acenaphthylene             | 14.30 | 152  | 551146   | 20.081 | ng/uL | 99       |
| 49) 3-Nitroaniline             | 14.47 | 138  | 87072    | 21.383 | ng/uL | 98       |
| 50) Acenaphthene               | 14.64 | 153  | 389958   | 19.734 | ng/uL | 99       |
| 51) 2,4-Dinitrophenol          | 14.68 | 184  | 31205    | 16.307 | ng/uL | 97       |
| 53) 4-Nitrophenol              | 14.78 | 109  | 50418    | 19.034 | ng/uL | 96       |
| 54) Dibenzofuran               | 14.98 | 168  | 548165   | 20.341 | ng/uL | 99       |
| 55) 2,4-Dinitrotoluene         | 14.93 | 165  | 136705   | 22.207 | ng/uL | 90       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.20 | 232  | 138583   | 23.345 | ng/uL | 94       |
| 57) Diethylphthalate           | 15.40 | 149  | 443840   | 19.823 | ng/uL | 98       |
| 59) Fluorene                   | 15.63 | 166  | 457568   | 20.023 | ng/uL | 98       |
| 60) 4-Chlorophenyl-phenylether | 15.62 | 204  | 244314   | 20.349 | ng/uL | 95       |
| 61) 4-Nitroaniline             | 15.63 | 138  | 94875    | 20.900 | ng/uL | 93       |
| 64) 4,6-Dinitro-2-methylphenol | 15.70 | 198  | 67920    | 19.382 | ng/uL | 94       |
| 65) N-Nitrosodiphenylamine     | 15.83 | 169  | 394860   | 20.659 | ng/uL | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.51 | 248  | 171712   | 21.021 | ng/uL | 94       |
| 67) Hexachlorobenzene          | 16.64 | 284  | 197864   | 21.458 | ng/uL | 96       |
| 68) Atrazine                   | 16.78 | 200  | 163964   | 21.459 | ng/uL | 100      |
| 69) Pentachlorophenol          | 16.97 | 266  | 110256   | 22.434 | ng/uL | 96       |
| 70) Phenanthrene               | 17.36 | 178  | 733232   | 19.986 | ng/uL | 100      |
| 72) Anthracene                 | 17.45 | 178  | 742743   | 20.319 | ng/uL | 98       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.38 | 216  | 211095   | 21.385 | ng/uL | 99       |
| 74) Pentachlorobenzene         | 14.90 | 250  | 218169   | 21.885 | ng/uL | 96       |
| 75) Carbazole                  | 17.72 | 167  | 645546   | 22.008 | ng/uL | 99       |
| 76) Di-n-butylphthalate        | 18.29 | 149  | 761996   | 21.113 | ng/uL | 99       |
| 77) Fluoranthene               | 19.37 | 202  | 889068   | 22.093 | ng/uL | 99       |
| 80) Pyrene                     | 19.73 | 202  | 877731   | 20.665 | ng/uL | 98       |
| 81) Butylbenzylphthalate       | 20.63 | 149  | 327207   | 22.444 | ng/uL | 93       |
| 82) 3,3'-Dichlorobenzidine     | 21.47 | 252  | 302416   | 22.431 | ng/uL | 96       |
| 83) Benzo(a)anthracene         | 21.55 | 228  | 846400   | 20.460 | ng/uL | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.48 | 149  | 463930   | 21.875 | ng/uL | 99       |
| 85) Chrysene                   | 21.61 | 228  | 817047   | 20.318 | ng/uL | 100      |
| 87) Di-n-octyl phthalate       | 22.65 | 149  | 808492   | 22.796 | ng/uL | 100      |
| 88) Benzo(b)fluoranthene       | 23.69 | 252  | 904382   | 19.723 | ng/uL | 100      |
| 89) Benzo(k)fluoranthene       | 23.76 | 252  | 898948   | 19.967 | ng/uL | 99       |
| 91) Benzo(a)pyrene             | 24.53 | 252  | 834637   | 19.849 | ng/uL | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 28.19 | 276  | 1060209  | 20.189 | ng/uL | 98       |
| 93) Dibenzo(a,h)anthracene     | 28.25 | 278  | 878124   | 20.440 | ng/uL | 98       |
| 94) Benzo(g,h,i)perylene       | 29.28 | 276  | 888438m  | 20.016 | ng/uL |          |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG080620\  
Quantitation Report (QT Reviewed)

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Acq On : 6 Aug 2020 12:00  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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Quant Time: Aug 06 12:42:56 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG072320MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 06 10:37:54 2020  
Response via : Initial Calibration

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

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

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```
Acq On       : 6 Aug 2020 12:00
Operator    : CG/JU
Sample      : SSTDCCC020
Misc        :
ALS Vial     : 2      Sample Multipl
```

**Instrument :**  
BNA\_G  
**ClientSampleId :**  
SSTD02040

**Manual Integrations**  
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