

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\  
 Data File : BP000165.D  
 Acq On : 10 Sep 2019 15:21  
 Operator : HP/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleID :  
 SSTD02015

Manual Integrations  
 APPROVED

mohammad  
 9/12/2019 9:29:08 AM

Quant Time: Sep 11 02:06:35 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP090519MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 10 02:43:18 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.89  | 152  | 417337   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 8.18  | 136  | 1614692  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 9.95  | 164  | 893450   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 11.45 | 188  | 1616883  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 14.15 | 240  | 1201356  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 15.69 | 264  | 1256037  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 2.61  | 96  | 83605   | 7.10  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.51  | 99  | 715216  | 20.12 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.58  | 67  | 372672  | 20.05 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.66  | 132 | 594386  | 20.68 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 7.26  | 113 | 556398  | 20.84 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 7.45  | 128 | 268608  | 21.91 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 7.78  | 143 | 273717  | 22.93 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 8.02  | 165 | 536611  | 21.18 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 8.23  | 131 | 711110  | 22.70 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 9.63  | 166 | 1372033 | 20.80 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 9.79  | 160 | 1817150 | 22.29 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 10.03 | 143 | 236651  | 20.07 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 10.47 | 176 | 1144356 | 20.43 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 10.53 | 200 | 146192  | 19.67 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 11.51 | 188 | 1634777 | 20.88 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 12.90 | 212 | 1600419 | 22.40 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 15.59 | 264 | 1308909 | 20.10 | ng/ul | 0.00 |

## Target Compounds

|                                |      |     |         |        | Ovalue    |
|--------------------------------|------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 2.68 | 88  | 83484   | 6.987  | ng/uL 98  |
| 4) Benzaldehyde                | 6.44 | 77  | 453169  | 25.231 | ng/ul 98  |
| 6) Phenol                      | 6.52 | 94  | 725609  | 19.586 | ng/ul 99  |
| 8) Bis(2-Chloroethyl)ether     | 6.62 | 93  | 585251  | 20.169 | ng/ul 95  |
| 10) 2-Chlorophenol             | 6.68 | 128 | 599963  | 19.770 | ng/ul 99  |
| 11) 2-Methylphenol             | 7.13 | 108 | 567366  | 20.478 | ng/ul 100 |
| 12) 2,2'-oxybis(1-Chloropropan | 7.17 | 45  | 559578  | 20.121 | ng/ul 97  |
| 14) Acetophenone               | 7.30 | 105 | 841473  | 21.002 | ng/ul 97  |
| 15) N-Nitroso-di-n-propylamine | 7.30 | 70  | 367713  | 19.539 | ng/ul 97  |
| 16) 4-Methylphenol             | 7.29 | 108 | 566249  | 20.556 | ng/ul 99  |
| 17) Hexachloroethane           | 7.41 | 117 | 239988  | 20.012 | ng/ul 97  |
| 20) Nitrobenzene               | 7.47 | 77  | 591898  | 20.444 | ng/ul 96  |
| 21) Isophorone                 | 7.71 | 82  | 1128072 | 19.910 | ng/ul 99  |
| 23) 2-Nitrophenol              | 7.79 | 139 | 295127  | 21.398 | ng/ul# 90 |
| 24) 2,4-Dimethylphenol         | 7.82 | 107 | 623144  | 20.220 | ng/ul 98  |
| 25) Bis(2-Chloroethoxy)methane | 7.92 | 93  | 751946  | 20.111 | ng/ul 99  |
| 27) 2,4-Dichlorophenol         | 8.03 | 162 | 516300  | 20.313 | ng/ul 98  |
| 28) Naphthalene                | 8.20 | 128 | 1786086 | 20.307 | ng/ul 100 |
| 30) 4-Chloroaniline            | 8.24 | 127 | 705756  | 22.213 | ng/ul 99  |
| 31) Hexachlorobutadiene        | 8.33 | 225 | 331785  | 20.773 | ng/ul 99  |
| 32) Caprolactam                | 8.58 | 113 | 214397m | 24.597 | ng/ul     |
| 33) 4-Chloro-3-methylphenol    | 8.72 | 107 | 539456  | 20.930 | ng/ul 100 |
| 34) 2-Methylnaphthalene        | 8.89 | 142 | 1218089 | 20.279 | ng/ul 99  |

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 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD02015

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 10 02:43:18 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 8.99  | 142  | 1183008  | 20.638 | ng/ul  | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 9.06  | 216  | 588013   | 19.774 | ng/ul  | 97       |
| 38) Hexachlorocyclopentadiene  | 9.06  | 237  | 289908   | 18.468 | ng/ul  | 99       |
| 39) 2,4,6-Trichlorophenol      | 9.17  | 196  | 370300   | 21.002 | ng/ul  | 100      |
| 40) 2,4,5-Trichlorophenol      | 9.21  | 196  | 381832   | 19.967 | ng/ul  | 98       |
| 41) 1,1'-Biphenyl              | 9.36  | 154  | 1456653  | 19.794 | ng/ul  | 99       |
| 42) 2-Chloronaphthalene        | 9.39  | 162  | 1155344  | 20.166 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 9.48  | 65   | 271267   | 21.104 | ng/ul  | 93       |
| 45) Dimethylphthalate          | 9.66  | 163  | 1335962  | 20.068 | ng/ul  | 100      |
| 46) 2,6-Dinitrotoluene         | 9.72  | 165  | 272195   | 21.765 | ng/ul  | 98       |
| 48) Acenaphthylene             | 9.80  | 152  | 1673393  | 19.631 | ng/ul  | 100      |
| 49) 3-Nitroaniline             | 9.89  | 138  | 302574   | 22.248 | ng/ul  | 98       |
| 50) Acenaphthene               | 9.98  | 153  | 1175530  | 19.917 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 9.99  | 184  | 94616    | 17.838 | ng/ul# | 1        |
| 53) 4-Nitrophenol              | 10.04 | 109  | 172137   | 19.620 | ng/ul  | 96       |
| 54) Dibenzofuran               | 10.15 | 168  | 1604774  | 19.818 | ng/ul  | 99       |
| 55) 2,4-Dinitrotoluene         | 10.13 | 165  | 368977   | 21.280 | ng/ul  | 96       |
| 56) 2,3,4,6-Tetrachlorophenol  | 10.27 | 232  | 306434   | 21.033 | ng/ul  | 97       |
| 57) Diethylphthalate           | 10.37 | 149  | 1337905  | 20.503 | ng/ul  | 100      |
| 59) Fluorene                   | 10.50 | 166  | 1231471  | 19.660 | ng/ul  | 99       |
| 60) 4-Chlorophenyl-phenylether | 10.49 | 204  | 623652   | 19.732 | ng/ul  | 95       |
| 61) 4-Nitroaniline             | 10.50 | 138  | 291749   | 22.285 | ng/ul  | 99       |
| 64) 4,6-Dinitro-2-methylphenol | 10.54 | 198  | 153056   | 18.913 | ng/ul  | 94       |
| 65) N-Nitrosodiphenylamine     | 10.61 | 169  | 1094982  | 20.166 | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 10.99 | 248  | 396038   | 20.979 | ng/ul# | 91       |
| 67) Hexachlorobenzene          | 11.06 | 284  | 407760   | 20.417 | ng/ul# | 89       |
| 68) Atrazine                   | 11.15 | 200  | 463319   | 24.154 | ng/ul  | 99       |
| 69) Pentachlorophenol          | 11.25 | 266  | 226828   | 20.264 | ng/ul  | 100      |
| 70) Phenanthrene               | 11.47 | 178  | 1887657  | 19.742 | ng/ul  | 100      |
| 72) Anthracene                 | 11.53 | 178  | 1886171  | 20.047 | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 9.35  | 216  | 571633   | 19.949 | ng/uL  | 99       |
| 74) Pentachlorobenzene         | 10.12 | 250  | 538679   | 20.755 | ng/uL  | 98       |
| 75) Carbazole                  | 11.68 | 167  | 1667303  | 20.726 | ng/ul  | 100      |
| 76) Di-n-butylphthalate        | 12.03 | 149  | 2168057  | 22.253 | ng/ul  | 100      |
| 77) Fluoranthene               | 12.69 | 202  | 2025669  | 21.004 | ng/ul  | 98       |
| 80) Pvrene                     | 12.92 | 202  | 1960310  | 21.628 | ng/ul# | 92       |
| 81) Butylbenzylphthalate       | 13.56 | 149  | 921293   | 25.904 | ng/ul  | 100      |
| 82) 3,3'-Dichlorobenzidine     | 14.10 | 252  | 655227   | 22.322 | ng/ul  | 98       |
| 83) Benzo(a)anthracene         | 14.13 | 228  | 1822168  | 20.953 | ng/ul  | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 14.16 | 149  | 1294504  | 26.210 | ng/ul  | 99       |
| 85) Chrysene                   | 14.17 | 228  | 1572736  | 19.708 | ng/ul  | 100      |
| 87) Di-n-octyl phthalate       | 14.79 | 149  | 2196526  | 27.236 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 15.23 | 252  | 1664069  | 19.989 | ng/ul  | 99       |
| 89) Benzo(k)fluoranthene       | 15.26 | 252  | 1593136  | 20.302 | ng/ul  | 99       |
| 91) Benzo(a)pyrene             | 15.62 | 252  | 1550334  | 19.849 | ng/ul  | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 17.26 | 276  | 1756163  | 19.544 | ng/ul  | 99       |
| 93) Dibenzo(a,h)anthracene     | 17.28 | 278  | 1452499  | 19.516 | ng/ul  | 99       |
| 94) Benzo(a,h,i)perylene       | 17.73 | 276  | 1385800  | 18.497 | ng/ul  | 99       |

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Data File : BP000165.D  
Acq On : 10 Sep 2019 15:21  
Operator : HP/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD02015

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Quant Title : SVOA CALIBRATION  
QLast Update : Tue Sep 10 02:43:18 2019  
Response via : Initial Calibration

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

