

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP030520\  
 Data File : BP002086.D  
 Acq On : 06 Mar 2020 00:27  
 Operator : CG/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD02099

Quant Time: Mar 06 07:58:41 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP022720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 06 07:37:34 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.64  | 152  | 305066   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.41 | 136  | 1242319  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.27 | 164  | 775019   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.03 | 188  | 1659606  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.12 | 240  | 1574415  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.33 | 264  | 1804875  | 20.00 | ng/ul | -0.01    |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.16  | 96  | 54849   | 7.73  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 6.82  | 99  | 453474  | 19.49 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.97  | 67  | 243474  | 19.76 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.17  | 132 | 389075  | 19.33 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.35  | 113 | 359895  | 18.82 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.78  | 128 | 175652  | 18.41 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 188768  | 17.58 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.03 | 165 | 399476  | 19.31 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 10.55 | 131 | 470228  | 20.74 | ng/ul | 0.00  |
| 44) Dimethylphthalate-d6       | 13.69 | 166 | 1107888 | 18.89 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 13.96 | 160 | 1315041 | 18.75 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 14.47 | 143 | 174237  | 15.54 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 15.27 | 176 | 964007  | 18.99 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.39 | 200 | 135063  | 14.76 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 17.12 | 188 | 1455971 | 19.52 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 19.37 | 212 | 1664468 | 20.20 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 23.19 | 264 | 1807164 | 19.57 | ng/ul | -0.01 |

## Target Compounds

|                                |       |     |         |        | Ovalue    |
|--------------------------------|-------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.20  | 88  | 54452   | 7.196  | ng/uL 96  |
| 4) Benzaldehyde                | 6.78  | 77  | 221780  | 16.219 | ng/ul 100 |
| 6) Phenol                      | 6.85  | 94  | 484523  | 20.056 | ng/ul 97  |
| 8) Bis(2-Chloroethyl)ether     | 7.07  | 93  | 373863  | 19.722 | ng/ul 99  |
| 10) 2-Chlorophenol             | 7.20  | 128 | 414059  | 20.063 | ng/ul 99  |
| 11) 2-Methylphenol             | 8.08  | 108 | 368041  | 19.484 | ng/ul 100 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.17  | 45  | 405574  | 20.042 | ng/ul 98  |
| 14) Acetophenone               | 8.45  | 105 | 566700  | 20.096 | ng/ul 98  |
| 15) N-Nitroso-di-n-propylamine | 8.44  | 70  | 261582  | 19.464 | ng/ul 97  |
| 16) 4-Methylphenol             | 8.40  | 108 | 404247  | 20.034 | ng/ul 99  |
| 17) Hexachloroethane           | 8.71  | 117 | 158092  | 19.280 | ng/ul 95  |
| 20) Nitrobenzene               | 8.82  | 77  | 398968  | 18.763 | ng/ul 98  |
| 21) Isophorone                 | 9.35  | 82  | 746712  | 18.508 | ng/ul 100 |
| 23) 2-Nitrophenol              | 9.53  | 139 | 215763  | 18.473 | ng/ul 98  |
| 24) 2,4-Dimethylphenol         | 9.60  | 107 | 421771  | 19.371 | ng/ul 99  |
| 25) Bis(2-Chloroethoxy)methane | 9.83  | 93  | 491423  | 19.235 | ng/ul 99  |
| 27) 2,4-Dichlorophenol         | 10.06 | 162 | 406086  | 19.750 | ng/ul 99  |
| 28) Naphthalene                | 10.46 | 128 | 1285210 | 19.310 | ng/ul 100 |
| 30) 4-Chloroaniline            | 10.57 | 127 | 494466  | 21.942 | ng/ul 98  |
| 31) Hexachlorobutadiene        | 10.76 | 225 | 275433  | 19.421 | ng/ul 99  |
| 32) Caprolactam                | 11.30 | 113 | 106606  | 15.784 | ng/ul 95  |
| 33) 4-Chloro-3-methylphenol    | 11.70 | 107 | 378394  | 18.519 | ng/ul 99  |
| 34) 2-Methylnaphthalene        | 12.08 | 142 | 933666  | 19.974 | ng/ul 100 |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP030520\  
 Data File : BP002086.D  
 Acq On : 06 Mar 2020 00:27  
 Operator : CG/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD02099

Quant Time: Mar 06 07:58:41 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP022720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 06 07:37:34 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.30 | 142  | 892376   | 19.226 | ng/ul | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.46 | 216  | 529901   | 20.510 | ng/ul | 99       |
| 38) Hexachlorocyclopentadiene  | 12.45 | 237  | 295742   | 19.530 | ng/ul | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.70 | 196  | 315933   | 18.589 | ng/ul | 99       |
| 40) 2,4,5-Trichlorophenol      | 12.77 | 196  | 347559   | 19.390 | ng/ul | 100      |
| 41) 1,1'-Biphenyl              | 13.10 | 154  | 1222933  | 20.285 | ng/ul | 100      |
| 42) 2-Chloronaphthalene        | 13.14 | 162  | 937668   | 19.675 | ng/ul | 99       |
| 43) 2-Nitroaniline             | 13.34 | 65   | 199010   | 17.536 | ng/ul | 99       |
| 45) Dimethylphthalate          | 13.73 | 163  | 1121835  | 18.529 | ng/ul | 100      |
| 46) 2,6-Dinitrotoluene         | 13.85 | 165  | 220286   | 17.929 | ng/ul | 98       |
| 48) Acenaphthylene             | 13.99 | 152  | 1421070  | 19.106 | ng/ul | 100      |
| 49) 3-Nitroaniline             | 14.17 | 138  | 211514   | 18.602 | ng/ul | 94       |
| 50) Acenaphthene               | 14.34 | 153  | 971286   | 19.423 | ng/ul | 98       |
| 51) 2,4-Dinitrophenol          | 14.39 | 184  | 144029   | 20.527 | ng/ul | 98       |
| 53) 4-Nitrophenol              | 14.49 | 109  | 131534   | 16.243 | ng/ul | 97       |
| 54) Dibenzofuran               | 14.67 | 168  | 1416477  | 20.054 | ng/ul | 99       |
| 55) 2,4-Dinitrotoluene         | 14.64 | 165  | 321119   | 18.247 | ng/ul | 100      |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.90 | 232  | 294829   | 17.956 | ng/ul | 98       |
| 57) Diethylphthalate           | 15.12 | 149  | 1099816  | 18.404 | ng/ul | 99       |
| 59) Fluorene                   | 15.33 | 166  | 1094376  | 19.389 | ng/ul | 98       |
| 60) 4-Chlorophenyl-phenylether | 15.33 | 204  | 574395   | 19.312 | ng/ul | 98       |
| 61) 4-Nitroaniline             | 15.34 | 138  | 215120   | 18.060 | ng/ul | 99       |
| 64) 4,6-Dinitro-2-methylphenol | 15.41 | 198  | 155774   | 15.632 | ng/ul | 96       |
| 65) N-Nitrosodiphenylamine     | 15.54 | 169  | 966011   | 20.446 | ng/ul | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.22 | 248  | 374756   | 19.704 | ng/ul | 95       |
| 67) Hexachlorobenzene          | 16.34 | 284  | 412557   | 19.103 | ng/ul | 99       |
| 68) Atrazine                   | 16.50 | 200  | 341877   | 18.323 | ng/ul | 99       |
| 69) Pentachlorophenol          | 16.68 | 266  | 210656   | 15.796 | ng/ul | 99       |
| 70) Phenanthrene               | 17.07 | 178  | 1769720  | 19.356 | ng/ul | 99       |
| 72) Anthracene                 | 17.16 | 178  | 1774227  | 19.599 | ng/ul | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.07 | 216  | 525769   | 19.720 | ng/uL | 99       |
| 74) Pentachlorobenzene         | 14.60 | 250  | 505386   | 19.196 | ng/uL | 99       |
| 75) Carbazole                  | 17.43 | 167  | 1585615  | 20.801 | ng/ul | 100      |
| 76) Di-n-butylphthalate        | 18.01 | 149  | 1750596  | 17.916 | ng/ul | 100      |
| 77) Fluoranthene               | 19.04 | 202  | 2087057  | 19.788 | ng/ul | 100      |
| 80) Pvrene                     | 19.40 | 202  | 2153344  | 20.088 | ng/ul | 99       |
| 81) Butylbenzylphthalate       | 20.29 | 149  | 790163   | 18.168 | ng/ul | 98       |
| 82) 3,3'-Dichlorobenzidine     | 21.04 | 252  | 639087   | 17.778 | ng/ul | 99       |
| 83) Benzo(a)anthracene         | 21.10 | 228  | 2197755  | 20.633 | ng/ul | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.06 | 149  | 1194753  | 18.718 | ng/ul | 100      |
| 85) Chrysene                   | 21.16 | 228  | 2024219  | 19.919 | ng/ul | 99       |
| 87) Di-n-octyl phthalate       | 21.92 | 149  | 1940954  | 18.233 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.66 | 252  | 2218128  | 19.597 | ng/ul | 99       |
| 89) Benzo(k)fluoranthene       | 22.71 | 252  | 2206836  | 19.918 | ng/ul | 100      |
| 91) Benzo(a)pyrene             | 23.23 | 252  | 2128506  | 20.715 | ng/ul | 100      |
| 92) Indeno(1,2,3-cd)pyrene     | 25.55 | 276  | 2543537  | 19.099 | ng/ul | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.56 | 278  | 2138302  | 19.419 | ng/ul | 99       |
| 94) Benzo(a,h,i)perylene       | 26.23 | 276  | 2131369  | 18.959 | ng/ul | 98       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP030520\  
Data File : BP002086.D  
Acq On : 06 Mar 2020 00:27  
Operator : CG/JU  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD02099

Quant Time: Mar 06 07:58:41 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP022720MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Mar 06 07:37:34 2020  
Response via : Initial Calibration

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

