

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\  
 Data File : BP001865.D  
 Acq On : 24 Feb 2020 17:44  
 Operator : CG/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleID :  
 SSTD02083

Manual Integrations  
 APPROVED

mohammad  
 2/25/2020 3:42:21 PM

Quant Time: Feb 25 05:54:25 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP021820MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 25 05:05:58 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.65  | 152  | 382647   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 1627256  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.29 | 164  | 1049702  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.04 | 188  | 2348376  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.13 | 240  | 2153896  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.35 | 264  | 2400584  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.17  | 96  | 61513   | 6.68  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.83  | 99  | 567714  | 19.64 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.99  | 67  | 305729  | 18.20 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.19  | 132 | 494804  | 21.18 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.36  | 113 | 483189  | 21.07 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 246954  | 22.42 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.51  | 143 | 278406  | 27.95 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.05 | 165 | 537733  | 22.70 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.56 | 131 | 603923  | 21.55 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.70 | 166 | 1582709 | 21.49 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.97 | 160 | 1934022 | 21.32 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.49 | 143 | 286935  | 22.37 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.29 | 176 | 1380798 | 20.84 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 234514  | 25.33 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.14 | 188 | 2104654 | 20.34 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.38 | 212 | 2381355 | 21.54 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.21 | 264 | 2501742 | 21.58 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue   |
|--------------------------------|-------|-----|---------|--------|----------|
| 2) 1,4-Dioxane                 | 3.20  | 88  | 65084   | 6.587  | ng/uL 99 |
| 4) Benzaldehyde                | 6.79  | 77  | 369789  | 20.894 | ng/ul 97 |
| 6) Phenol                      | 6.86  | 94  | 591066  | 19.291 | ng/ul 98 |
| 8) Bis(2-Chloroethyl)ether     | 7.08  | 93  | 468640  | 18.967 | ng/ul 97 |
| 10) 2-Chlorophenol             | 7.22  | 128 | 510891  | 20.669 | ng/ul 98 |
| 11) 2-Methylphenol             | 8.09  | 108 | 475735  | 20.508 | ng/ul 99 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.19  | 45  | 514037  | 17.461 | ng/ul 96 |
| 14) Acetophenone               | 8.46  | 105 | 731952  | 20.655 | ng/ul 97 |
| 15) N-Nitroso-di-n-propylamine | 8.46  | 70  | 346224  | 21.072 | ng/ul 97 |
| 16) 4-Methylphenol             | 8.42  | 108 | 516112  | 20.664 | ng/ul 96 |
| 17) Hexachloroethane           | 8.72  | 117 | 204549  | 20.803 | ng/ul 97 |
| 20) Nitrobenzene               | 8.83  | 77  | 544286  | 19.933 | ng/ul 94 |
| 21) Isophorone                 | 9.36  | 82  | 1053325 | 20.988 | ng/ul 98 |
| 23) 2-Nitrophenol              | 9.55  | 139 | 300184  | 25.148 | ng/ul 97 |
| 24) 2,4-Dimethylphenol         | 9.62  | 107 | 562114  | 20.967 | ng/ul 98 |
| 25) Bis(2-Chloroethoxy)methane | 9.85  | 93  | 657295  | 19.353 | ng/ul 99 |
| 27) 2,4-Dichlorophenol         | 10.08 | 162 | 528752  | 21.747 | ng/ul 99 |
| 28) Naphthalene                | 10.48 | 128 | 1699319 | 19.542 | ng/ul 99 |
| 30) 4-Chloroaniline            | 10.58 | 127 | 601515  | 21.242 | ng/ul 99 |
| 31) Hexachlorobutadiene        | 10.78 | 225 | 365388  | 21.253 | ng/ul 99 |
| 32) Caprolactam                | 11.32 | 113 | 170373m | 23.194 | ng/ul    |
| 33) 4-Chloro-3-methylphenol    | 11.72 | 107 | 533627  | 23.051 | ng/ul 98 |
| 34) 2-Methylnaphthalene        | 12.09 | 142 | 1221275 | 20.374 | ng/ul 97 |

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD02083

Manual Integrations  
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Quant Time: Feb 25 05:54:25 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP021820MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 25 05:05:58 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.32 | 142  | 1211051  | 20.500 | ng/ul | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.47 | 216  | 693860   | 20.294 | ng/ul | 99       |
| 38) Hexachlorocyclopentadiene  | 12.46 | 237  | 403213   | 21.833 | ng/ul | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.71 | 196  | 446910   | 24.455 | ng/ul | 99       |
| 40) 2,4,5-Trichlorophenol      | 12.78 | 196  | 479985   | 23.856 | ng/ul | 99       |
| 41) 1,1'-Biphenyl              | 13.12 | 154  | 1639801  | 20.038 | ng/ul | 99       |
| 42) 2-Chloronaphthalene        | 13.15 | 162  | 1293898  | 20.025 | ng/ul | 100      |
| 43) 2-Nitroaniline             | 13.35 | 65   | 305311   | 23.540 | ng/ul | 92       |
| 45) Dimethylphthalate          | 13.75 | 163  | 1628382  | 21.003 | ng/ul | 100      |
| 46) 2,6-Dinitrotoluene         | 13.86 | 165  | 343267   | 25.405 | ng/ul | 90       |
| 48) Acenaphthylene             | 14.00 | 152  | 2022554  | 20.553 | ng/ul | 99       |
| 49) 3-Nitroaniline             | 14.19 | 138  | 309345   | 22.488 | ng/ul | 95       |
| 50) Acenaphthene               | 14.35 | 153  | 1342222  | 19.961 | ng/ul | 99       |
| 51) 2,4-Dinitrophenol          | 14.39 | 184  | 148001   | 26.396 | ng/ul | 95       |
| 53) 4-Nitrophenol              | 14.50 | 109  | 210756   | 22.192 | ng/ul | 91       |
| 54) Dibenzofuran               | 14.69 | 168  | 1917647  | 20.250 | ng/ul | 99       |
| 55) 2,4-Dinitrotoluene         | 14.65 | 165  | 487649   | 24.512 | ng/ul | 93       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.92 | 232  | 424810   | 26.124 | ng/ul | 96       |
| 57) Diethylphthalate           | 15.13 | 149  | 1618121  | 21.827 | ng/ul | 98       |
| 59) Fluorene                   | 15.34 | 166  | 1520865  | 20.188 | ng/ul | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.35 | 204  | 797432   | 20.397 | ng/ul | 99       |
| 61) 4-Nitroaniline             | 15.35 | 138  | 347688   | 22.386 | ng/ul | 95       |
| 64) 4,6-Dinitro-2-methylphenol | 15.42 | 198  | 253016   | 24.504 | ng/ul | 98       |
| 65) N-Nitrosodiphenylamine     | 15.56 | 169  | 1330502  | 19.761 | ng/ul | 100      |
| 66) 4-Bromophenyl-phenylether  | 16.24 | 248  | 534139   | 20.723 | ng/ul | 99       |
| 67) Hexachlorobenzene          | 16.36 | 284  | 609403   | 20.487 | ng/ul | 97       |
| 68) Atrazine                   | 16.52 | 200  | 540352   | 22.260 | ng/ul | 99       |
| 69) Pentachlorophenol          | 16.70 | 266  | 344993   | 23.956 | ng/ul | 99       |
| 70) Phenanthrene               | 17.08 | 178  | 2524520  | 19.493 | ng/ul | 99       |
| 72) Anthracene                 | 17.17 | 178  | 2536191  | 19.825 | ng/ul | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.08 | 216  | 753314   | 19.568 | ng/uL | 98       |
| 74) Pentachlorobenzene         | 14.62 | 250  | 740136   | 19.724 | ng/uL | 99       |
| 75) Carbazole                  | 17.45 | 167  | 2234026  | 20.940 | ng/ul | 99       |
| 76) Di-n-butylphthalate        | 18.02 | 149  | 2747547  | 23.402 | ng/ul | 99       |
| 77) Fluoranthene               | 19.06 | 202  | 3081578  | 22.220 | ng/ul | 96       |
| 80) Pvrene                     | 19.41 | 202  | 3083004  | 21.043 | ng/ul | 98       |
| 81) Butylbenzylphthalate       | 20.30 | 149  | 1286438  | 28.283 | ng/ul | 95       |
| 82) 3,3'-Dichlorobenzidine     | 21.06 | 252  | 1088424  | 23.367 | ng/ul | 99       |
| 83) Benzo(a)anthracene         | 21.12 | 228  | 2999972  | 21.150 | ng/ul | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.07 | 149  | 1867793  | 24.785 | ng/ul | 99       |
| 85) Chrysene                   | 21.17 | 228  | 2833736  | 20.450 | ng/ul | 99       |
| 87) Di-n-octyl phthalate       | 21.94 | 149  | 3230304  | 26.869 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.69 | 252  | 3227753  | 22.851 | ng/ul | 99       |
| 89) Benzo(k)fluoranthene       | 22.73 | 252  | 2888955  | 19.731 | ng/ul | 99       |
| 91) Benzo(a)pyrene             | 23.25 | 252  | 2739114  | 20.691 | ng/ul | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.58 | 276  | 3117025  | 18.850 | ng/ul | 98       |
| 93) Dibenzo(a,h)anthracene     | 25.60 | 278  | 2619415  | 18.860 | ng/ul | 98       |
| 94) Benzo(a,h,i)perylene       | 26.26 | 276  | 2550136  | 18.141 | ng/ul | 97       |

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Data File : BP001865.D  
Acq On : 24 Feb 2020 17:44  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION  
QLast Update : Tue Feb 25 05:05:58 2020  
Response via : Initial Calibration

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

