

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
 Data File : BN014287.D  
 Acq On : 15 Apr 2021 16:50  
 Operator : CG/JU  
 Sample : PB135733BS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SLCS733

Manual Integrations  
 APPROVED

mohammad  
 4/16/2021 11:30:57 AM

Quant Time: Apr 16 01:37:02 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  | 143749   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.47 | 136  | 567571   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.33 | 164  | 336549   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.08 | 188  | 664563   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.28 | 240  | 707792   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.51 | 264  | 732269   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 24313   | 6.44  | ng/uL | 0.00  |
| 4) Pyridine-d5                 | 3.62  | 84  | 320002  | 31.98 | ng/ul | 0.00  |
| 7) Phenol-d5                   | 6.86  | 99  | 416688  | 34.67 | ng/ul | -0.01 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.03  | 67  | 241505  | 33.56 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4          | 7.23  | 132 | 339590  | 35.93 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8          | 8.39  | 113 | 317767  | 34.88 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5            | 8.84  | 128 | 149205  | 36.38 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4           | 9.56  | 143 | 154749  | 38.53 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3      | 10.09 | 165 | 317114  | 36.65 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4         | 10.60 | 131 | 382766  | 29.82 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6       | 13.75 | 166 | 885454  | 36.32 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8          | 14.02 | 160 | 1146149 | 39.08 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4           | 14.53 | 143 | 171418  | 36.25 | ng/ul | 0.00  |
| 60) Fluorene-d10               | 15.33 | 176 | 770459  | 35.75 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylphenol | 15.45 | 200 | 131061  | 36.85 | ng/ul | 0.00  |
| 73) Anthracene-d10             | 17.18 | 188 | 1220002 | 39.33 | ng/ul | 0.00  |
| 81) Pyrene-d10                 | 19.48 | 212 | 1363646 | 39.69 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12         | 23.37 | 264 | 1484249 | 38.21 | ng/ul | 0.00  |

## Target Compounds

|                                |       |     |         |        | Ovalue    |
|--------------------------------|-------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.25  | 88  | 48613   | 12.635 | ng/uL# 95 |
| 5) Pyridine                    | 3.63  | 79  | 335220  | 32.599 | ng/ul 96  |
| 6) Benzaldehyde                | 6.84  | 77  | 219196  | 36.126 | ng/ul 98  |
| 8) Phenol                      | 6.89  | 94  | 438224  | 33.924 | ng/ul 99  |
| 10) Bis(2-Chloroethyl)ether    | 7.12  | 93  | 340006  | 33.467 | ng/ul 99  |
| 12) 2-Chlorophenol             | 7.26  | 128 | 350644  | 34.696 | ng/ul 97  |
| 13) 2-Methylphenol             | 8.13  | 108 | 319149  | 34.071 | ng/ul 95  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.22  | 45  | 372405  | 33.160 | ng/ul 99  |
| 16) Acetophenone               | 8.51  | 105 | 515186  | 33.647 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.50  | 70  | 249411  | 35.035 | ng/ul 99  |
| 18) 4-Methylphenol             | 8.46  | 108 | 347473  | 34.252 | ng/ul 97  |
| 19) Hexachloroethane           | 8.76  | 117 | 146697  | 34.532 | ng/ul 97  |
| 22) Nitrobenzene               | 8.88  | 77  | 382073  | 35.216 | ng/ul 99  |
| 23) Isophorone                 | 9.40  | 82  | 662682  | 36.318 | ng/ul 99  |
| 25) 2-Nitrophenol              | 9.59  | 139 | 177786  | 37.462 | ng/ul 99  |
| 26) 2,4-Dimethylphenol         | 9.65  | 107 | 372512  | 35.259 | ng/ul 98  |
| 27) Bis(2-Chloroethoxy)methane | 9.89  | 93  | 428212  | 34.447 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 10.12 | 162 | 319080  | 35.936 | ng/ul 98  |
| 30) Naphthalene                | 10.52 | 128 | 1076587 | 34.224 | ng/ul 100 |
| 32) 4-Chloroaniline            | 10.63 | 127 | 386109  | 29.234 | ng/ul 99  |
| 33) Hexachlorobutadiene        | 10.81 | 225 | 203075  | 35.030 | ng/ul 97  |
| 34) Caprolactam                | 11.39 | 113 | 92570m  | 37.404 | ng/ul     |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
 Data File : BN014287.D  
 Acq On : 15 Apr 2021 16:50  
 Operator : CG/JU  
 Sample : PB135733BS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SLCS733

Manual Integrations  
 APPROVED

mohammad  
 4/16/2021 11:30:57 AM

Quant Time: Apr 16 01:37:02 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  | 321607   | 36.645 | ng/ul  | 98       |
| 36) 2-Methylnaphthalene        | 12.14 | 142  | 745821   | 34.796 | ng/ul  | 99       |
| 37) 1-Methylnaphthalene        | 12.36 | 142  | 719005   | 33.687 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.51 | 216  | 388533   | 34.108 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene  | 12.50 | 237  | 224175   | 31.691 | ng/ul  | 97       |
| 41) 2,4,6-Trichlorophenol      | 12.75 | 196  | 237119   | 36.016 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol      | 12.82 | 196  | 259689   | 35.543 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl              | 13.16 | 154  | 923914   | 33.379 | ng/ul  | 97       |
| 44) 2-Chloronaphthalene        | 13.20 | 162  | 727103   | 33.604 | ng/ul  | 98       |
| 45) 2-Nitroaniline             | 13.40 | 65   | 191883   | 38.253 | ng/ul  | 93       |
| 47) Dimethylphthalate          | 13.79 | 163  | 915535   | 36.172 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene         | 13.91 | 165  | 177450   | 39.977 | ng/ul  | 98       |
| 50) Acenaphthylene             | 14.05 | 152  | 1076563  | 34.497 | ng/ul  | 99       |
| 51) 3-Nitroaniline             | 14.23 | 138  | 175161   | 32.026 | ng/ul  | 95       |
| 52) Acenaphthene               | 14.40 | 153  | 780787   | 34.300 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol          | 14.44 | 184  | 84710    | 32.694 | ng/ul  | 95       |
| 55) 4-Nitrophenol              | 14.55 | 109  | 143183   | 35.338 | ng/ul  | 95       |
| 56) Dibenzofuran               | 14.73 | 168  | 1083137  | 34.191 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene         | 14.70 | 165  | 265040   | 41.333 | ng/ul  | 92       |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.96 | 232  | 211619   | 36.815 | ng/ul  | 97       |
| 59) Diethylphthalate           | 15.17 | 149  | 933248   | 36.046 | ng/ul  | 99       |
| 61) Fluorene                   | 15.39 | 166  | 917515   | 35.729 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenylether | 15.38 | 204  | 450981   | 35.662 | ng/ul  | 98       |
| 63) 4-Nitroaniline             | 15.40 | 138  | 200045   | 37.576 | ng/ul  | 100      |
| 66) 4,6-Dinitro-2-methylphenol | 15.46 | 198  | 139208   | 36.964 | ng/ul# | 94       |
| 67) N-Nitrosodiphenylamine     | 15.59 | 169  | 768154   | 37.149 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether  | 16.27 | 248  | 265653   | 37.785 | ng/ul  | 97       |
| 69) Hexachlorobenzene          | 16.39 | 284  | 318529   | 37.955 | ng/ul  | 95       |
| 70) Atrazine                   | 16.55 | 200  | 274002   | 37.287 | ng/ul  | 99       |
| 71) Pentachlorophenol          | 16.73 | 266  | 175051   | 35.875 | ng/ul  | 91       |
| 72) Phenanthrene               | 17.12 | 178  | 1434389  | 36.968 | ng/ul  | 100      |
| 74) Anthracene                 | 17.22 | 178  | 1459858  | 37.042 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.13 | 216  | 378816   | 34.615 | ng/uL  | 97       |
| 76) Pentachlorobenzene         | 14.65 | 250  | 367542   | 33.791 | ng/uL  | 92       |
| 77) Carbazole                  | 17.48 | 167  | 1336379  | 37.484 | ng/ul  | 99       |
| 78) Di-n-butylphthalate        | 18.06 | 149  | 1567719  | 40.640 | ng/ul  | 99       |
| 80) Fluoranthene               | 19.15 | 202  | 1674538  | 38.569 | ng/ul  | 97       |
| 82) Pyrene                     | 19.51 | 202  | 1769064  | 38.266 | ng/ul  | 98       |
| 83) Butylbenzylphthalate       | 20.42 | 149  | 659225   | 43.430 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine     | 21.20 | 252  | 495776   | 39.165 | ng/ul  | 93       |
| 85) Benzo(a)anthracene         | 21.26 | 228  | 1793210  | 39.210 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.20 | 149  | 1074283  | 43.423 | ng/ul  | 99       |
| 87) Chrysene                   | 21.32 | 228  | 1751679  | 39.515 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate       | 22.09 | 149  | 1752777  | 38.338 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 22.84 | 252  | 1828519  | 37.325 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene       | 22.89 | 252  | 1860593  | 38.152 | ng/ul  | 99       |
| 93) Benzo(a)pyrene             | 23.42 | 252  | 1607054  | 37.078 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.76 | 276  | 2148838  | 37.406 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene     | 25.77 | 278  | 1818709  | 38.349 | ng/ul  | 95       |
| 96) Benzo(g,h,i)perylene       | 26.44 | 276  | 1734515  | 37.362 | ng/ul  | 99       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
Data File : BN014287.D  
Acq On : 15 Apr 2021 16:50  
Operator : CG/JU  
Sample : PB135733BS  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SLCS733

Manual Integrations  
APPROVED

mohammad  
4/16/2021 11:30:57 AM

Quant Time: Apr 16 01:37:02 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

