

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012021\  
 Data File : BN013435.D  
 Acq On : 20 Jan 2021 12:40  
 Operator : CG/JU  
 Sample : L5214-06  
 Misc : SFAM-MDL-ME-02  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 MDL-MED-SOIL-QT1-2021-02

Manual Integrations  
 APPROVED

mohammad  
 1/22/2021 11:50:57 AM

Quant Time: Jan 20 13:10:23 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN011521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 20 09:33:01 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.59  | 152  | 509925   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.35 | 136  | 2011879  | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.22 | 164  | 1105439  | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.97 | 188  | 2012981  | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.17 | 240  | 1576471  | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.35 | 264  | 1528087  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.19  | 96  | 42086   | 3.29  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.57  | 84  | 633558  | 19.08 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.76  | 99  | 1475297 | 35.97 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.93  | 67  | 901594  | 37.45 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.12  | 132 | 1298832 | 36.86 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.29  | 113 | 1201046 | 36.46 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.73  | 128 | 598725  | 38.43 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.44  | 143 | 613910  | 37.78 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 9.98  | 165 | 1117119 | 35.61 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.49 | 131 | 1609517 | 35.52 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.64 | 166 | 3015327 | 36.25 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 13.91 | 160 | 4067146 | 39.00 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.42 | 143 | 513670  | 32.26 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.22 | 176 | 2554527 | 34.97 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.33 | 200 | 376682  | 31.71 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.07 | 188 | 3572937 | 36.53 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.37 | 212 | 3425427 | 37.94 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.22 | 264 | 2950078 | 36.42 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue    |
|--------------------------------|-------|-----|--------|-------|-----------|
| 2) 1,4-Dioxane                 | 3.22  | 88  | 18207  | 1.378 | ng/uL 98  |
| 5) Pyridine                    | 3.59  | 79  | 135439 | 4.039 | ng/ul# 82 |
| 6) Benzaldehyde                | 6.74  | 77  | 82765  | 5.347 | ng/ul 97  |
| 8) Phenol                      | 6.79  | 94  | 175843 | 4.159 | ng/ul 96  |
| 10) Bis(2-Chloroethyl)ether    | 7.02  | 93  | 150017 | 4.330 | ng/ul 98  |
| 12) 2-Chlorophenol             | 7.15  | 128 | 151566 | 4.163 | ng/ul 97  |
| 13) 2-Methylphenol             | 8.02  | 108 | 116602 | 3.620 | ng/ul 98  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.12  | 45  | 206641 | 4.273 | ng/ul 98  |
| 16) Acetophenone               | 8.40  | 105 | 203772 | 4.077 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.39  | 70  | 85660  | 3.794 | ng/ul 98  |
| 18) 4-Methylphenol             | 8.35  | 108 | 139184 | 4.013 | ng/ul 98  |
| 19) Hexachloroethane           | 8.65  | 117 | 60535  | 4.051 | ng/ul 95  |
| 22) Nitrobenzene               | 8.77  | 77  | 149916 | 4.241 | ng/ul 97  |
| 23) Isophorone                 | 9.29  | 82  | 216637 | 3.458 | ng/ul 97  |
| 25) 2-Nitrophenol              | 9.47  | 139 | 77206  | 4.236 | ng/ul 99  |
| 26) 2,4-Dimethylphenol         | 9.54  | 107 | 143402 | 4.041 | ng/ul 98  |
| 27) Bis(2-Chloroethoxy)methane | 9.78  | 93  | 172668 | 4.004 | ng/ul 97  |
| 29) 2,4-Dichlorophenol         | 10.00 | 162 | 131230 | 4.180 | ng/ul 93  |
| 30) Naphthalene                | 10.40 | 128 | 486744 | 4.262 | ng/ul 98  |
| 32) 4-Chloroaniline            | 10.51 | 127 | 181197 | 4.146 | ng/ul 100 |
| 33) Hexachlorobutadiene        | 10.69 | 225 | 84480  | 4.119 | ng/ul 100 |
| 34) Caprolactam                | 11.29 | 113 | 21937m | 2.406 | ng/ul     |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012021\  
 Data File : BN013435.D  
 Acq On : 20 Jan 2021 12:40  
 Operator : CG/JU  
 Sample : L5214-06  
 Misc : SFAM-MDL-ME-02  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 MDL-MED-SOIL-QT1-2021-02

Manual Integrations  
 APPROVED

mohammad  
 1/22/2021 11:50:57 AM

Quant Time: Jan 20 13:10:23 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN011521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 20 09:33:01 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 4-Chloro-3-methylphenol    | 11.63 | 107  | 109130   | 3.487 | ng/ul  | 98       |
| 36) 2-Methylnaphthalene        | 12.02 | 142  | 320138   | 4.164 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene        | 12.24 | 142  | 309463   | 4.177 | ng/ul  | 97       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.39 | 216  | 156172   | 4.404 | ng/ul  | 100      |
| 40) Hexachlorocyclopentadiene  | 12.38 | 237  | 77070    | 3.521 | ng/ul# | 92       |
| 41) 2,4,6-Trichlorophenol      | 12.63 | 196  | 74329    | 3.454 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol      | 12.70 | 196  | 85186    | 3.635 | ng/ul  | 97       |
| 43) 1,1'-Biphenyl              | 13.05 | 154  | 399066   | 4.287 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene        | 13.08 | 162  | 312575   | 4.260 | ng/ul  | 99       |
| 45) 2-Nitroaniline             | 13.29 | 65   | 54002    | 3.275 | ng/ul  | 94       |
| 47) Dimethylphthalate          | 13.68 | 163  | 355084   | 4.140 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene         | 13.80 | 165  | 53132    | 3.233 | ng/ul# | 87       |
| 50) Acenaphthylene             | 13.94 | 152  | 453998   | 4.214 | ng/ul  | 100      |
| 51) 3-Nitroaniline             | 14.12 | 138  | 49013    | 3.213 | ng/ul  | 100      |
| 52) Acenaphthene               | 14.28 | 153  | 318958   | 4.150 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol          | 14.33 | 184  | 18239    | 1.982 | ng/ul# | 90       |
| 55) 4-Nitrophenol              | 14.43 | 109  | 43798    | 3.835 | ng/ul  | 92       |
| 56) Dibenzofuran               | 14.62 | 168  | 442566   | 4.249 | ng/ul  | 96       |
| 57) 2,4-Dinitrotoluene         | 14.59 | 165  | 80713    | 3.455 | ng/ul  | 94       |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.85 | 232  | 59742    | 3.160 | ng/ul  | 99       |
| 59) Diethylphthalate           | 15.06 | 149  | 307763   | 3.680 | ng/ul  | 99       |
| 61) Fluorene                   | 15.27 | 166  | 352978   | 4.251 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenylether | 15.27 | 204  | 170936   | 4.217 | ng/ul  | 96       |
| 63) 4-Nitroaniline             | 15.29 | 138  | 47461    | 3.497 | ng/ul  | 97       |
| 66) 4,6-Dinitro-2-methylphenol | 15.35 | 198  | 47740    | 3.706 | ng/ul# | 86       |
| 67) N-Nitrosodiphenylamine     | 15.49 | 169  | 278958   | 4.280 | ng/ul  | 95       |
| 68) 4-Bromophenyl-phenylether  | 16.17 | 248  | 91067    | 4.006 | ng/ul  | 93       |
| 69) Hexachlorobenzene          | 16.27 | 284  | 109912   | 4.276 | ng/ul  | 98       |
| 70) Atrazine                   | 16.45 | 200  | 71959    | 3.308 | ng/ul  | 98       |
| 71) Pentachlorophenol          | 16.62 | 266  | 35775    | 2.460 | ng/ul  | 98       |
| 72) Phenanthrene               | 17.01 | 178  | 513799   | 4.246 | ng/ul  | 99       |
| 74) Anthracene                 | 17.10 | 178  | 517612   | 4.242 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.00 | 216  | 148931   | 4.583 | ng/uL  | 99       |
| 76) Pentachlorobenzene         | 14.54 | 250  | 136063   | 4.302 | ng/uL  | 99       |
| 77) Carbazole                  | 17.37 | 167  | 393351   | 3.922 | ng/ul  | 99       |
| 78) Di-n-butylphthalate        | 17.96 | 149  | 360532   | 3.009 | ng/ul  | 98       |
| 80) Fluoranthene               | 19.03 | 202  | 475486   | 4.052 | ng/ul  | 100      |
| 82) Pyrene                     | 19.40 | 202  | 534022   | 4.416 | ng/ul  | 98       |
| 83) Butylbenzylphthalate       | 20.33 | 149  | 107302   | 2.527 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine     | 21.10 | 252  | 63523    | 2.150 | ng/ul  | 93       |
| 85) Benzo(a)anthracene         | 21.16 | 228  | 417285   | 3.975 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.12 | 149  | 158302   | 2.586 | ng/ul  | 99       |
| 87) Chrysene                   | 21.20 | 228  | 431013   | 4.167 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate       | 21.98 | 149  | 237442   | 2.380 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 22.70 | 252  | 404875   | 3.926 | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene       | 22.75 | 252  | 421778   | 4.158 | ng/ul  | 99       |
| 93) Benzo(a)pyrene             | 23.26 | 252  | 387382   | 4.203 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.53 | 276  | 473613   | 4.551 | ng/ul  | 95       |
| 95) Dibenzo(a,h)anthracene     | 25.54 | 278  | 405174   | 4.657 | ng/ul  | 97       |
| 96) Benzo(g,h,i)perylene       | 26.19 | 276  | 390915   | 4.477 | ng/ul  | 97       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN012021\  
Data File : BN013435.D  
Acq On : 20 Jan 2021 12:40  
Operator : CG/JU  
Sample : L5214-06  
Misc : SFAM-MDL-ME-02  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
MDL-MED-SOIL-QT1-2021-02

Manual Integrations  
APPROVED

mohammad  
1/22/2021 11:50:57 AM

Quant Time: Jan 20 13:10:23 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN011521.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Jan 20 09:33:01 2021  
Response via : Initial Calibration

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

|       |  |  |  |  |  |  |
|-------|--|--|--|--|--|--|
| ----- |  |  |  |  |  |  |
| ----- |  |  |  |  |  |  |

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

