

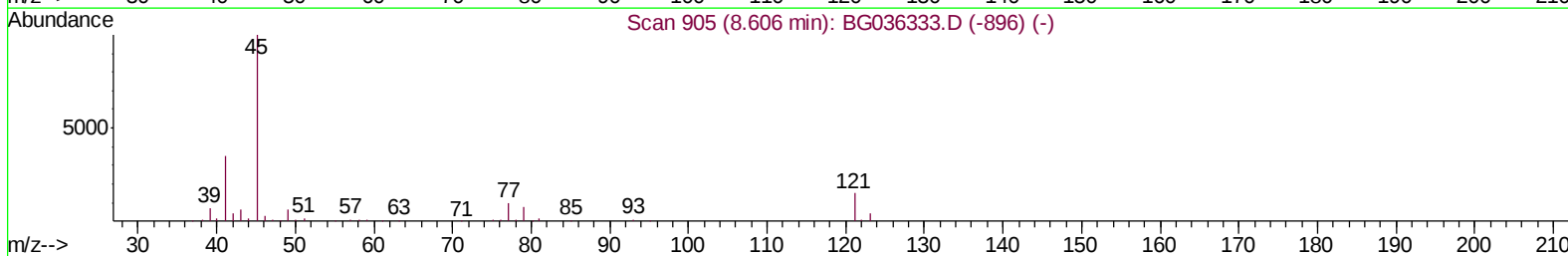
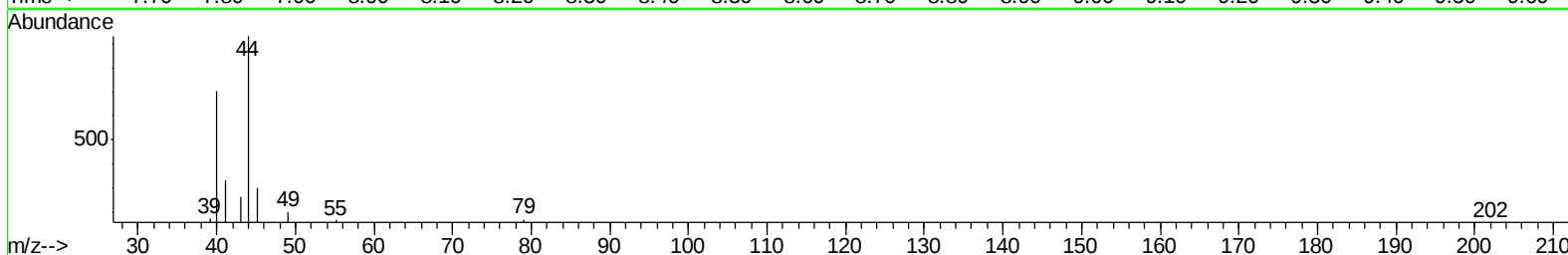
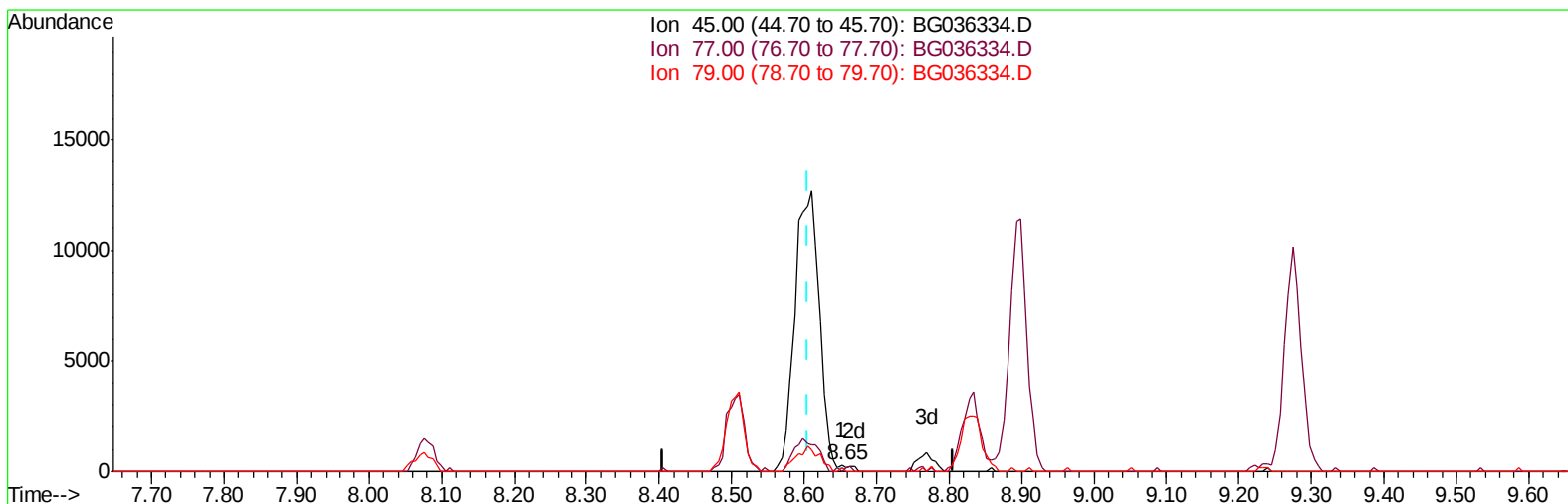
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036334.D  
 Acq On : 22 Aug 2018 9:27  
 Operator : JU/SJ  
 Sample : MDL-W-01  
 Misc : 1PPM/4PPM  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-W-01

Manual Integrations  
 APPROVED

Sohil  
 8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:27:44 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:22:46 2018  
 Response via : Initial Calibration



TIC: BG036334.D

(12) 2,2'-oxybis(1-Chloropropane)

8.652min (+0.046) 0.01ng/ul

response 75

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 45.00 | 100   | 100    |
| 77.00 | 15.60 | 0.00#  |
| 79.00 | 11.80 | 57.05# |
| 0.00  | 0.00  | 0.00   |

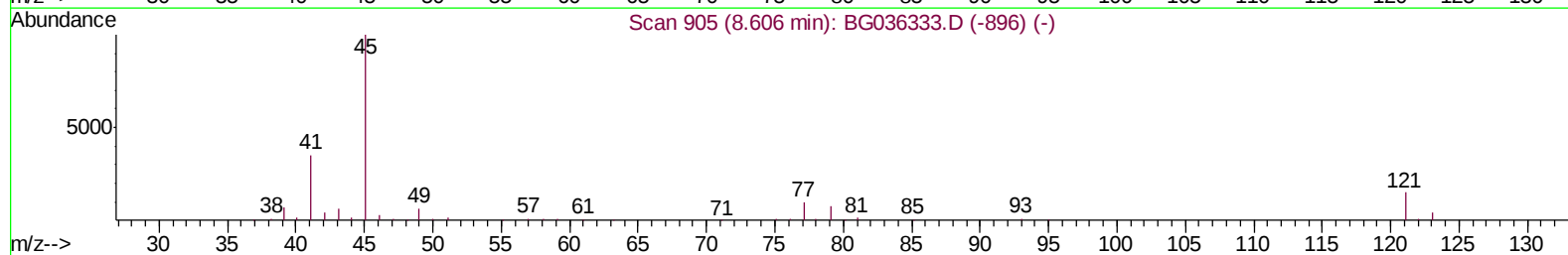
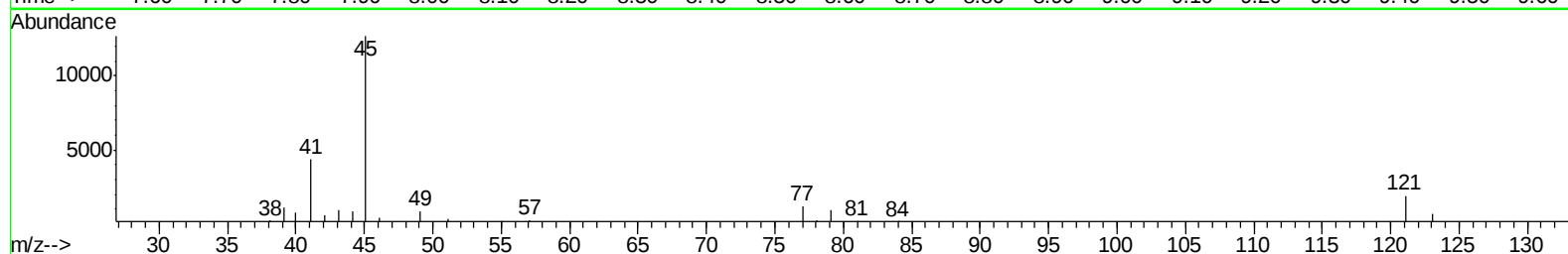
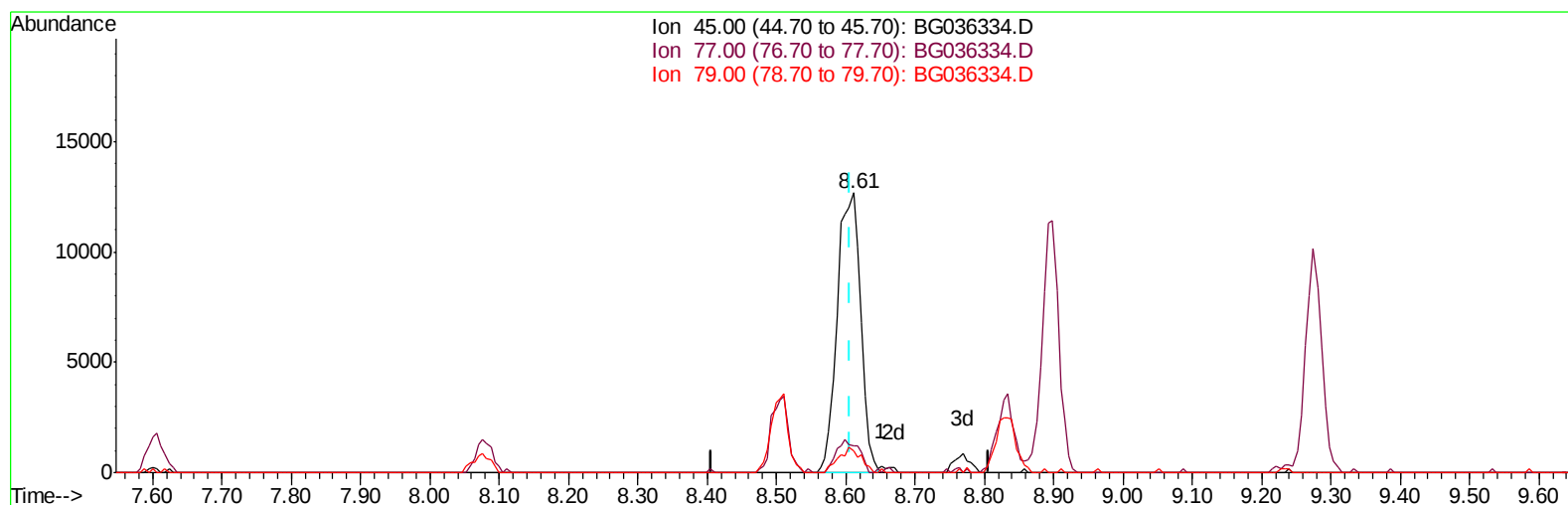
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036334.D  
 Acq On : 22 Aug 2018 9:27  
 Operator : JU/SJ  
 Sample : MDL-W-01  
 Misc : 1PPM/4PPM  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-W-01

Manual Integrations  
 APPROVED

Sohil  
 8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:27:44 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:22:46 2018  
 Response via : Initial Calibration



TIC: BG036334.D

(12) 2,2'-oxybis(1-Chloropropane)

8.611min (+0.004) 3.87ng/ul m

response 29731

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 45.00 | 100   | 100   |
| 77.00 | 15.60 | 9.61# |
| 79.00 | 11.80 | 7.77# |
| 0.00  | 0.00  | 0.00  |

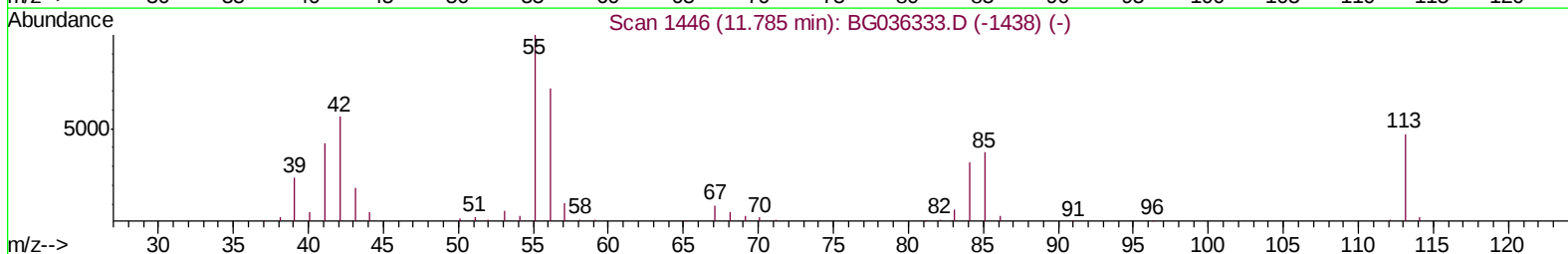
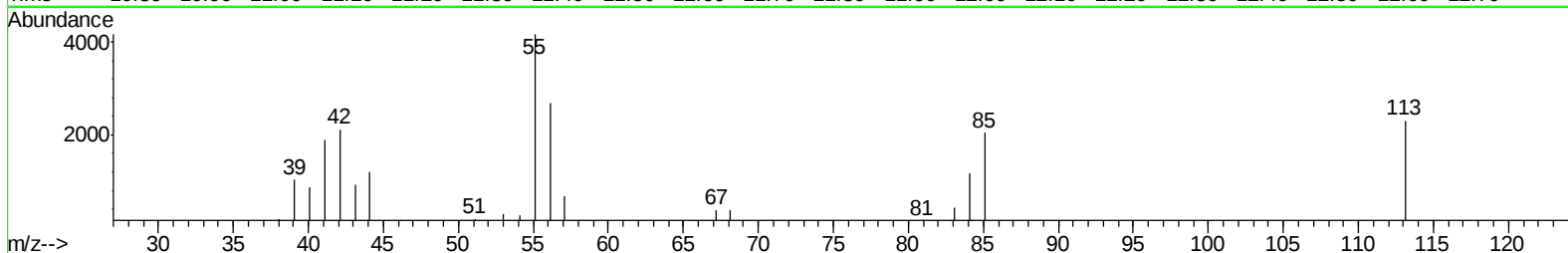
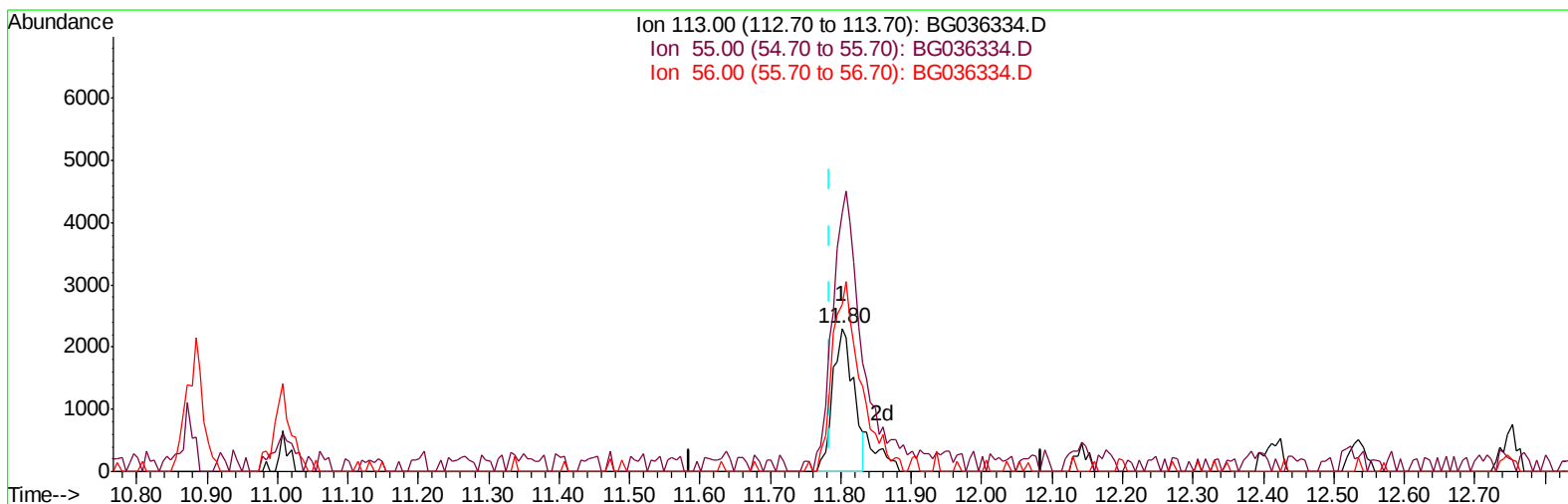
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036334.D  
Acq On : 22 Aug 2018 9:27  
Operator : JU/SJ  
Sample : MDL-W-01  
Misc : 1PPM/4PPM  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MDL-W-01

Manual Integrations  
APPROVED

Sohil  
8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:27:44 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:22:46 2018  
Response via : Initial Calibration



TIC: BG036334.D

(32) Caprolactam

11.801min (+0.016) 3.35ng/ul

response 4765

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 180.96 |
| 56.00  | 115.90 | 116.91 |
| 0.00   | 0.00   | 0.00   |

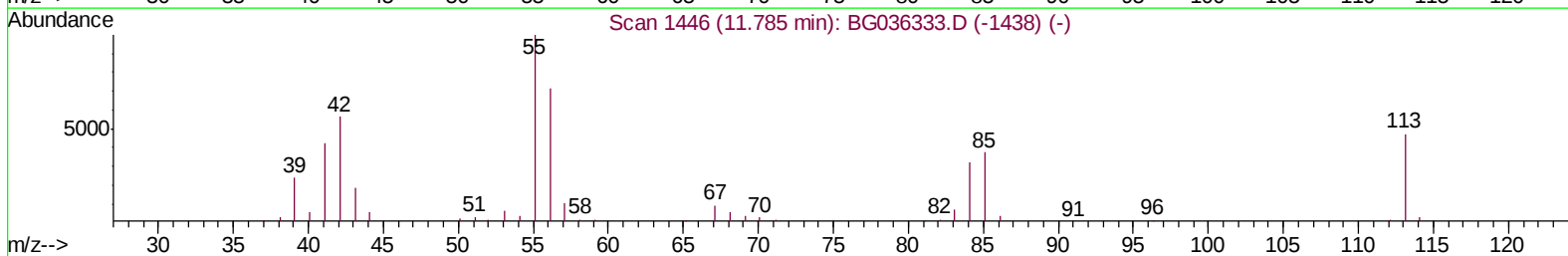
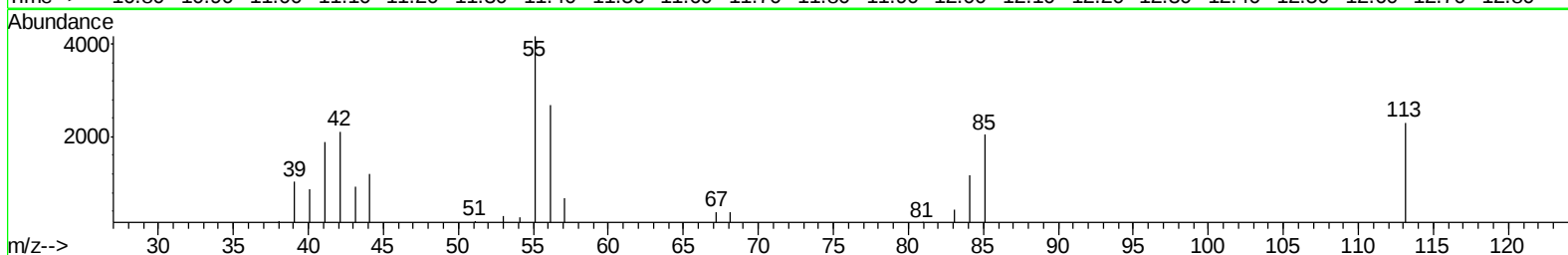
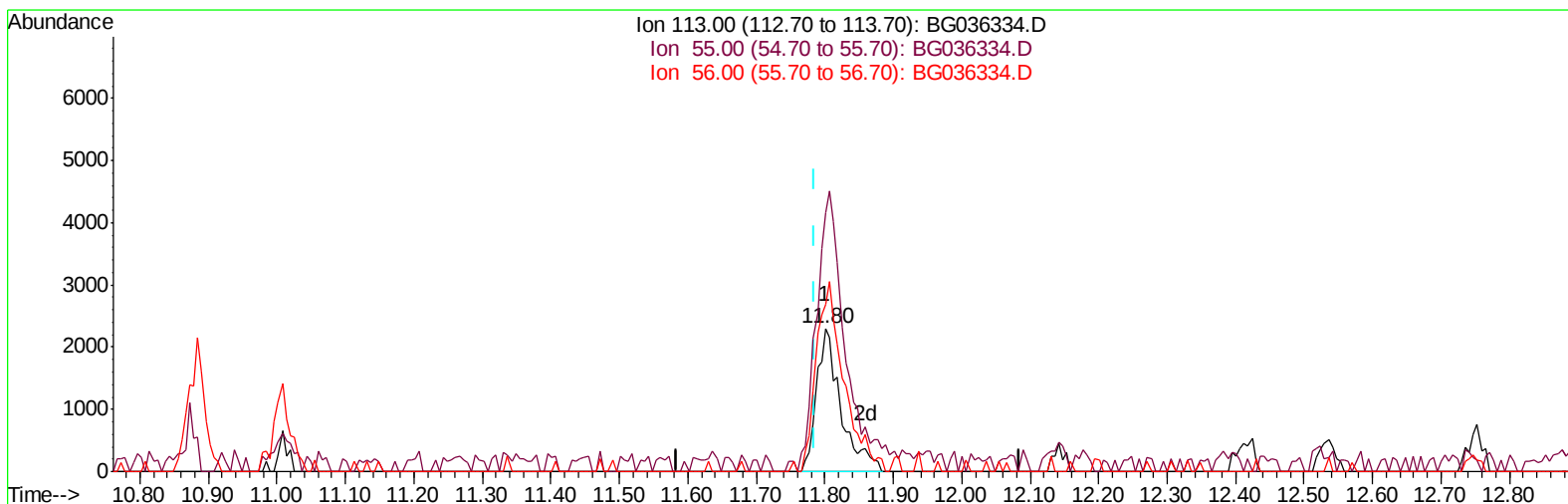
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036334.D  
Acq On : 22 Aug 2018 9:27  
Operator : JU/SJ  
Sample : MDL-W-01  
Misc : 1PPM/4PPM  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MDL-W-01

Quant Time: Aug 22 10:27:44 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:22:46 2018  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Sohil  
8/23/2018 10:58:39 AM



TIC: BG036334.D

(32) Caprolactam

11.801min (+0.016) 4.00ng/ul m

response 5682

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 180.96 |
| 56.00  | 115.90 | 116.91 |
| 0.00   | 0.00   | 0.00   |

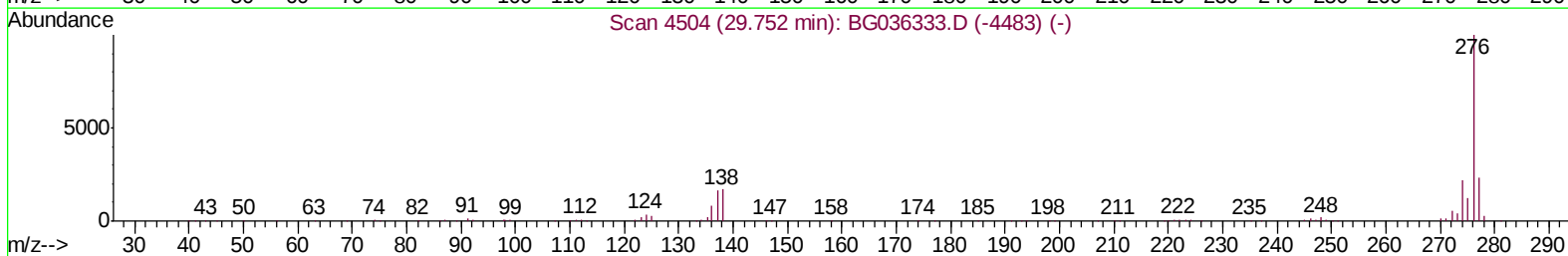
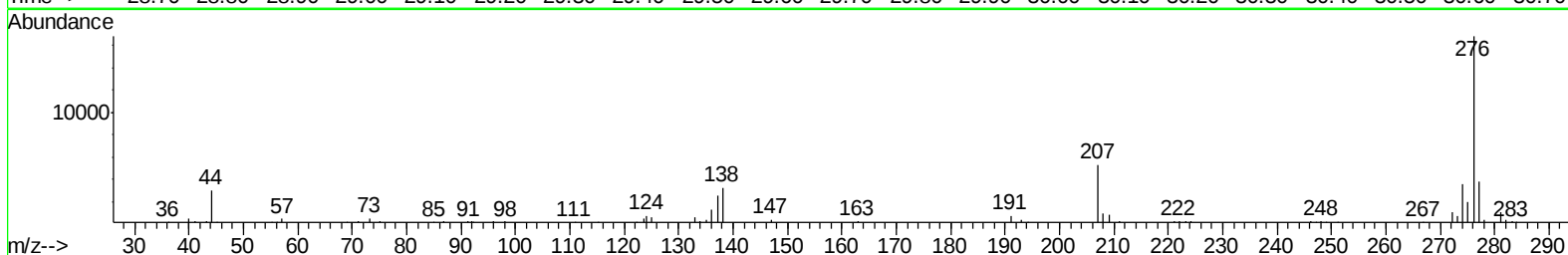
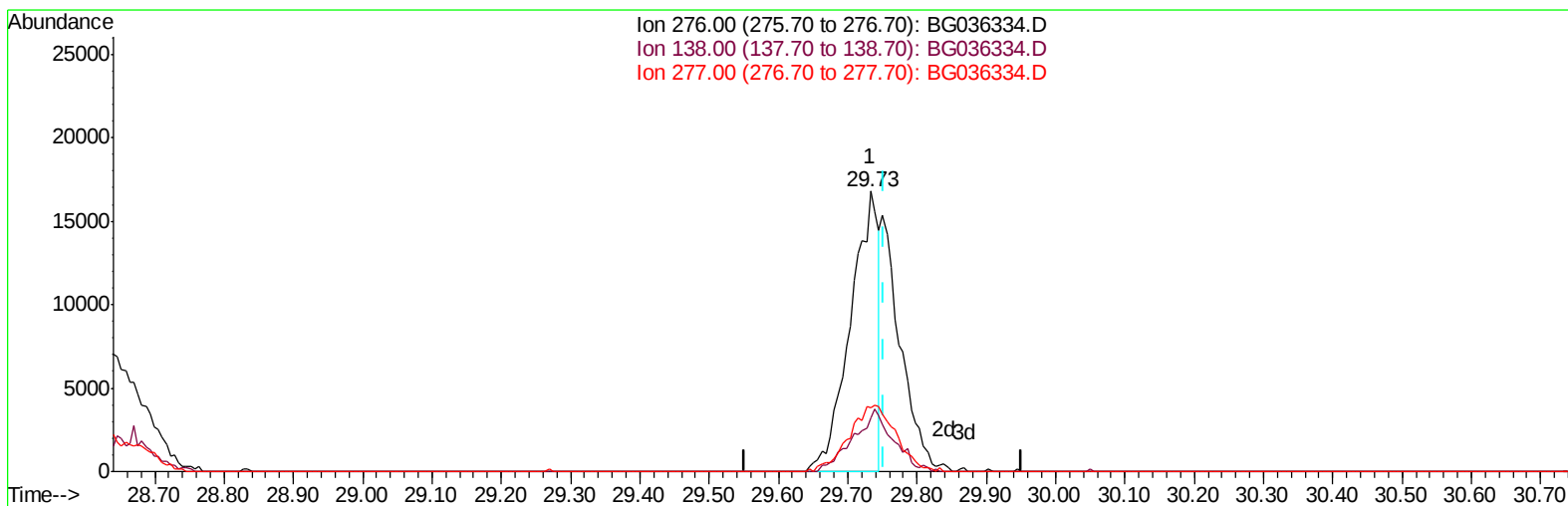
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036334.D  
Acq On : 22 Aug 2018 9:27  
Operator : JU/SJ  
Sample : MDL-W-01  
Misc : 1PPM/4PPM  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MDL-W-01

Manual Integrations  
APPROVED

Sohil  
8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:27:44 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:22:46 2018  
Response via : Initial Calibration



TIC: BG036334.D

(94) Benzo(g,h,i)perylene  
29.733min (-0.019) 2.37ng/ul  
response 47496

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 276.00 | 100   | 100    |
| 138.00 | 14.00 | 19.27# |
| 277.00 | 23.80 | 22.88  |
| 0.00   | 0.00  | 0.00   |

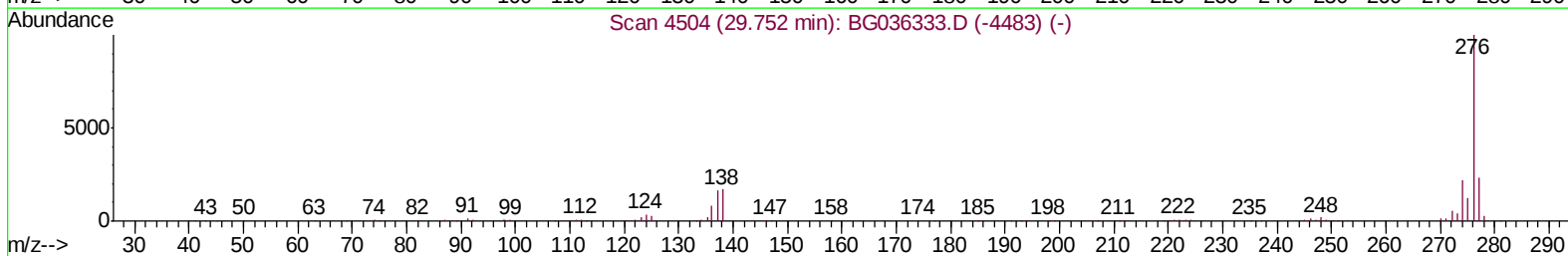
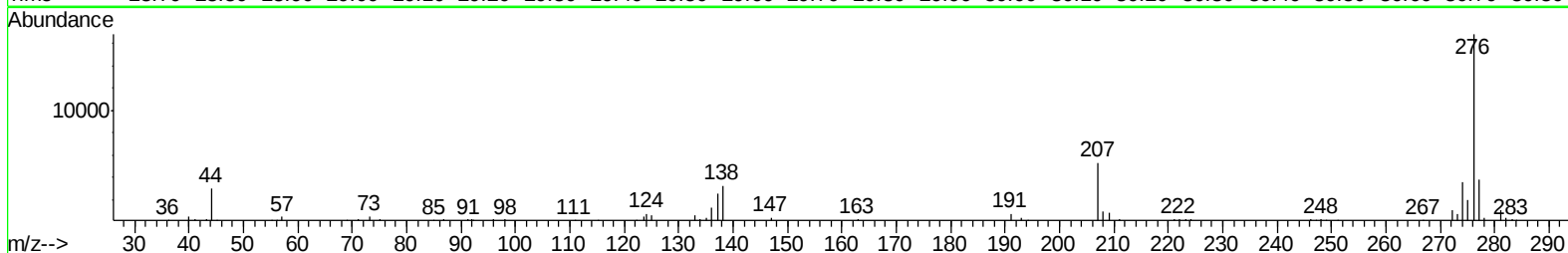
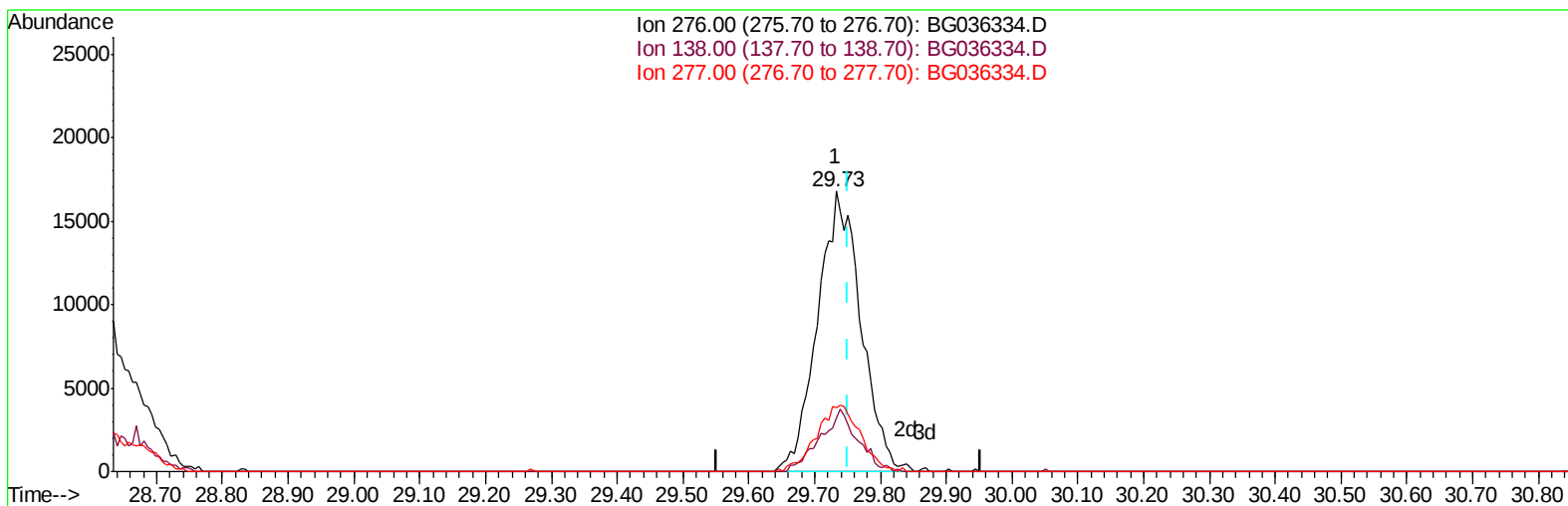
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036334.D  
Acq On : 22 Aug 2018 9:27  
Operator : JU/SJ  
Sample : MDL-W-01  
Misc : 1PPM/4PPM  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MDL-W-01

Manual Integrations  
APPROVED

Sohil  
8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:27:44 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:22:46 2018  
Response via : Initial Calibration



TIC: BG036334.D

(94) Benzo(g,h,i)perylene

29.733min (-0.019) 3.86ng/ul m

response 77377

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 276.00 | 100   | 100    |
| 138.00 | 14.00 | 19.27# |
| 277.00 | 23.80 | 22.88  |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036334.D  
 Acq On : 22 Aug 2018 9:27  
 Operator : JU/SJ  
 Sample : MDL-W-01  
 Misc : 1PPM/4PPM  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 MDL-W-01

Manual Integrations  
 APPROVED

Sohil  
 8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:32:31 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:22:46 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.08  | 152  | 50539    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.88 | 136  | 220535   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.70 | 164  | 138385   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.44 | 188  | 348903   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.71 | 240  | 375337   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.94 | 264  | 364728   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.52  | 96  | 7620   | 5.62  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.22  | 99  | 151069 | 32.13 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.40  | 67  | 105792 | 32.85 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.61  | 132 | 109711 | 32.32 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.77  | 113 | 123821 | 34.50 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.23  | 128 | 53222  | 39.56 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.96  | 143 | 52929  | 39.10 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.49 | 165 | 115269 | 31.82 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.01 | 131 | 160463 | 40.04 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.10 | 166 | 391702 | 33.85 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.39 | 160 | 469707 | 33.67 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.87 | 143 | 61662  | 36.03 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.69 | 176 | 329942 | 32.79 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.79 | 200 | 39705  | 31.42 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.54 | 188 | 536580 | 32.96 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.82 | 212 | 623321 | 32.41 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.72 | 264 | 598894 | 30.26 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |              | Qvalue |
|--------------------------------|-------|-----|--------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.56  | 88  | 2182   | 1.524 ng/uL# | 16     |
| 4) Benzaldehyde                | 7.22  | 77  | 7003   | 2.222 ng/ul# | 82     |
| 6) Phenol                      | 7.25  | 94  | 18567  | 3.951 ng/ul  | 96     |
| 8) Bis(2-Chloroethyl)ether     | 7.49  | 93  | 14819  | 4.247 ng/ul# | 83     |
| 10) 2-Chlorophenol             | 7.64  | 128 | 13179  | 3.881 ng/ul# | 88     |
| 11) 2-Methylphenol             | 8.50  | 108 | 13703  | 3.827 ng/ul  | 95     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.61  | 45  | 29731m | 3.870 ng/ul  |        |
| 14) Acetophenone               | 8.89  | 105 | 22865  | 4.156 ng/ul# | 85     |
| 15) N-Nitroso-di-n-propylamine | 8.88  | 70  | 12616  | 3.806 ng/ul# | 77     |
| 16) 4-Methylphenol             | 8.83  | 108 | 14233  | 3.800 ng/ul  | 91     |
| 17) Hexachloroethane           | 9.17  | 117 | 5734   | 4.238 ng/ul  | 85     |
| 20) Nitrobenzene               | 9.27  | 77  | 16376  | 4.290 ng/ul  | 96     |
| 21) Isophorone                 | 9.80  | 82  | 35859  | 3.870 ng/ul# | 96     |
| 23) 2-Nitrophenol              | 9.99  | 139 | 6670   | 4.655 ng/ul# | 93     |
| 24) 2,4-Dimethylphenol         | 10.04 | 107 | 16102  | 3.680 ng/ul  | 96     |
| 25) Bis(2-Chloroethoxy)methane | 10.29 | 93  | 19764  | 3.945 ng/ul  | 98     |
| 27) 2,4-Dichlorophenol         | 10.52 | 162 | 13528  | 3.941 ng/ul# | 84     |
| 28) Naphthalene                | 10.93 | 128 | 45096  | 3.990 ng/ul# | 93     |
| 30) 4-Chloroaniline            | 11.03 | 127 | 15132  | 3.875 ng/ul  | 91     |
| 31) Hexachlorobutadiene        | 11.22 | 225 | 9593   | 3.915 ng/ul  | 96     |
| 32) Caprolactam                | 11.80 | 113 | 5682m  | 4.000 ng/ul  |        |
| 33) 4-Chloro-3-methylphenol    | 12.14 | 107 | 15111  | 3.747 ng/ul  | 94     |
| 34) 2-Methylnaphthalene        | 12.54 | 142 | 34308  | 4.067 ng/ul  | 96     |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036334.D  
 Acq On : 22 Aug 2018 9:27  
 Operator : JU/SJ  
 Sample : MDL-W-01  
 Misc : 1PPM/4PPM  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-W-01

Manual Integrations  
 APPROVED

Sohil  
 8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:32:31 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:22:46 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.75 | 142  | 33143    | 4.028 | ng/ul# | 93       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.89 | 216  | 18710    | 3.969 | ng/ul  | 98       |
| 38) Hexachlorocyclopentadiene  | 12.88 | 237  | 10566    | 3.746 | ng/ul  | 99       |
| 39) 2,4,6-Trichlorophenol      | 13.13 | 196  | 10042    | 3.449 | ng/ul  | 92       |
| 40) 2,4,5-Trichlorophenol      | 13.19 | 196  | 12128    | 4.019 | ng/ul  | 91       |
| 41) 1,1'-Biphenyl              | 13.53 | 154  | 43284    | 3.996 | ng/ul# | 96       |
| 42) 2-Chloronaphthalene        | 13.58 | 162  | 35549    | 4.278 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 13.77 | 65   | 9108     | 4.189 | ng/ul# | 85       |
| 45) Dimethylphthalate          | 14.15 | 163  | 45488    | 4.097 | ng/ul  | 98       |
| 46) 2,6-Dinitrotoluene         | 14.27 | 165  | 6593     | 4.249 | ng/ul  | 96       |
| 48) Acenaphthylene             | 14.42 | 152  | 51794    | 3.938 | ng/ul# | 94       |
| 49) 3-Nitroaniline             | 14.59 | 138  | 6448     | 3.988 | ng/ul# | 92       |
| 50) Acenaphthene               | 14.76 | 153  | 37074    | 3.937 | ng/ul  | 98       |
| 51) 2,4-Dinitrophenol          | 14.79 | 184  | 2579     | 3.268 | ng/ul# | 88       |
| 53) 4-Nitrophenol              | 14.88 | 109  | 6163     | 3.880 | ng/ul# | 70       |
| 54) Dibenzofuran               | 15.10 | 168  | 54874    | 4.210 | ng/ul  | 98       |
| 55) 2,4-Dinitrotoluene         | 15.04 | 165  | 8239     | 3.839 | ng/ul# | 80       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.31 | 232  | 9852     | 3.486 | ng/ul# | 84       |
| 57) Diethylphthalate           | 15.51 | 149  | 46075    | 4.089 | ng/ul  | 97       |
| 59) Fluorene                   | 15.74 | 166  | 42286    | 4.024 | ng/ul  | 97       |
| 60) 4-Chlorophenyl-phenylether | 15.74 | 204  | 23067    | 4.029 | ng/ul  | 93       |
| 61) 4-Nitroaniline             | 15.74 | 138  | 5900     | 2.880 | ng/ul# | 84       |
| 64) 4,6-Dinitro-2-methylphenol | 15.80 | 198  | 5012     | 3.840 | ng/ul# | 64       |
| 65) N-Nitrosodiphenylamine     | 15.95 | 169  | 37833    | 3.871 | ng/ul  | 97       |
| 66) 4-Bromophenyl-phenylether  | 16.63 | 248  | 15197    | 3.938 | ng/ul  | 99       |
| 67) Hexachlorobenzene          | 16.75 | 284  | 15511    | 4.027 | ng/ul# | 95       |
| 68) Atrazine                   | 16.90 | 200  | 15422    | 3.825 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 17.08 | 266  | 7900     | 3.383 | ng/ul  | 90       |
| 70) Phenanthrene               | 17.48 | 178  | 73762    | 4.068 | ng/ul  | 96       |
| 72) Anthracene                 | 17.58 | 178  | 74175    | 4.065 | ng/ul  | 97       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.50 | 216  | 20191    | 3.787 | ng/uL  | 98       |
| 74) Pentachlorobenzene         | 15.01 | 250  | 18280    | 3.939 | ng/uL  | 94       |
| 75) Carbazole                  | 17.84 | 167  | 66587    | 4.341 | ng/ul  | 98       |
| 76) Di-n-butylphthalate        | 18.41 | 149  | 79741    | 3.979 | ng/ul  | 97       |
| 77) Fluoranthene               | 19.49 | 202  | 90484    | 4.304 | ng/ul# | 93       |
| 80) Pyrene                     | 19.85 | 202  | 89210    | 3.923 | ng/ul# | 93       |
| 81) Butylbenzylphthalate       | 20.74 | 149  | 33885    | 3.660 | ng/ul  | 95       |
| 82) 3,3'-Dichlorobenzidine     | 21.61 | 252  | 24998    | 3.233 | ng/ul  | 92       |
| 83) Benzo(a)anthracene         | 21.69 | 228  | 92660    | 3.907 | ng/ul  | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.62 | 149  | 49619    | 3.760 | ng/ul  | 97       |
| 85) Chrysene                   | 21.75 | 228  | 85074    | 3.911 | ng/ul  | 97       |
| 87) Di-n-octyl phthalate       | 22.85 | 149  | 80904    | 3.947 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.92 | 252  | 84838    | 3.928 | ng/ul# | 98       |
| 89) Benzo(k)fluoranthene       | 23.98 | 252  | 81457    | 3.912 | ng/ul  | 97       |
| 91) Benzo(a)pyrene             | 24.79 | 252  | 79408    | 3.886 | ng/ul  | 97       |
| 92) Indeno(1,2,3-cd)pyrene     | 28.59 | 276  | 94459    | 3.882 | ng/ul  | 98       |
| 93) Dibenzo(a,h)anthracene     | 28.68 | 278  | 75132    | 3.808 | ng/ul  | 97       |
| 94) Benzo(g,h,i)perylene       | 29.73 | 276  | 77377m   | 3.856 | ng/ul  |          |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036334.D  
Acq On : 22 Aug 2018 9:27  
Operator : JU/SJ  
Sample : MDL-W-01  
Misc : 1PPM/4PPM  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MDL-W-01

Manual Integrations  
APPROVED

Sohil  
8/23/2018 10:58:39 AM

Quant Time: Aug 22 10:32:31 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:22:46 2018  
Response via : Initial Calibration

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

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

