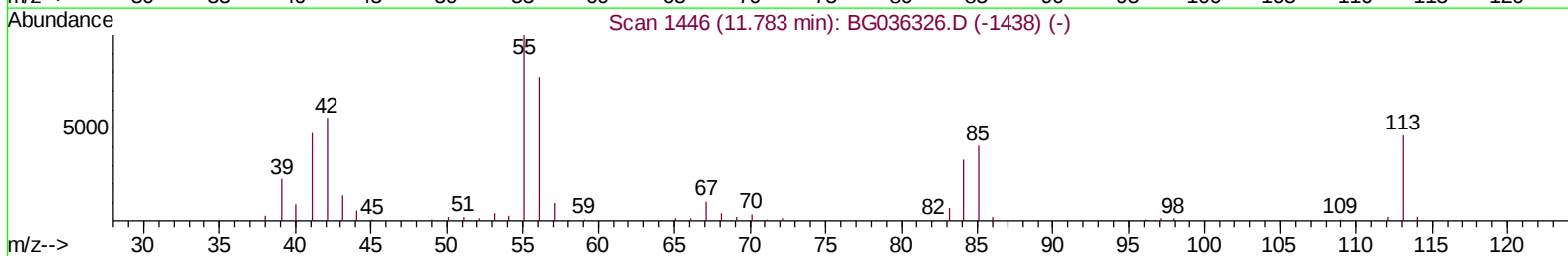
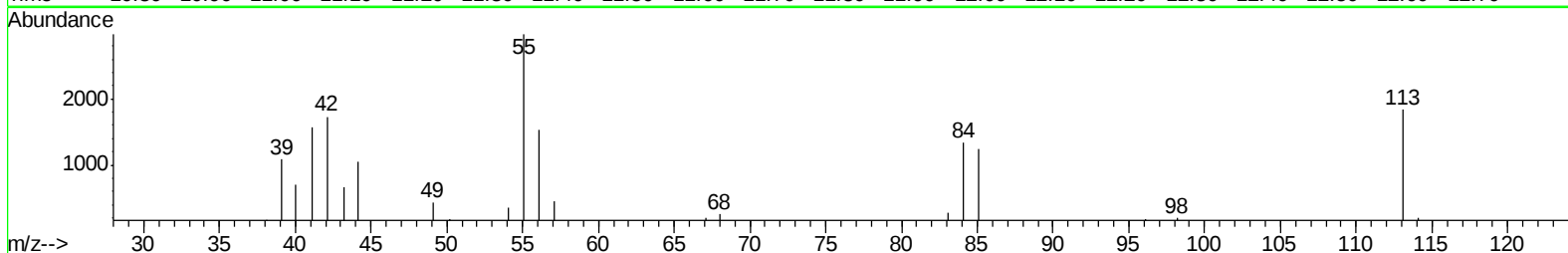
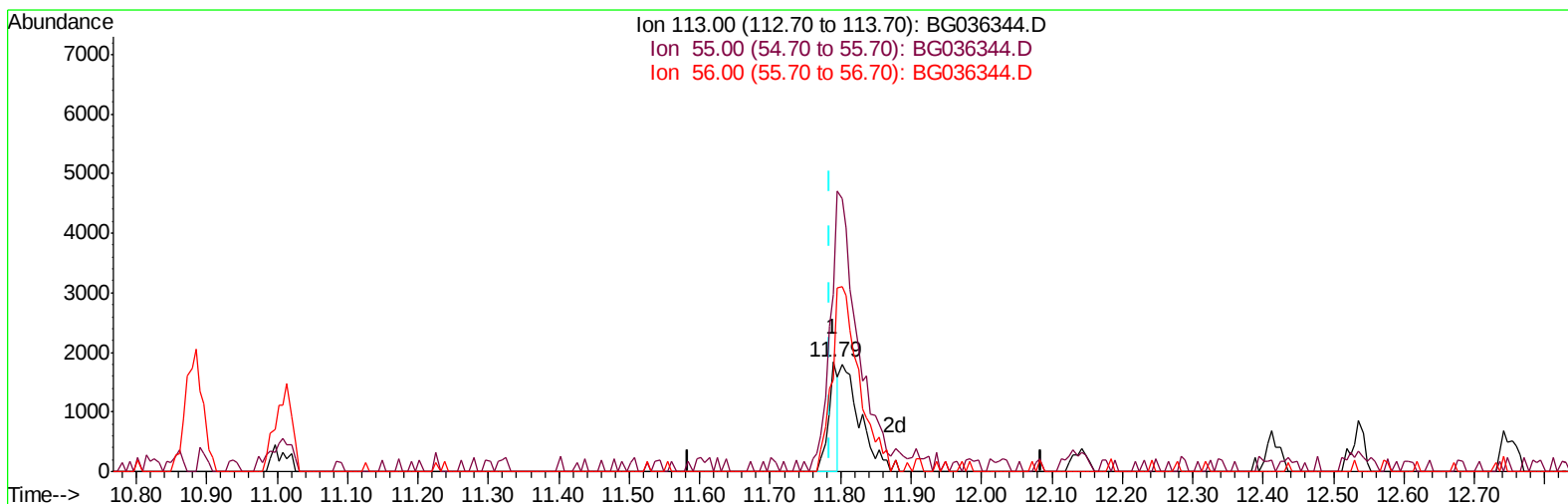


Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

(32) Caprolactam

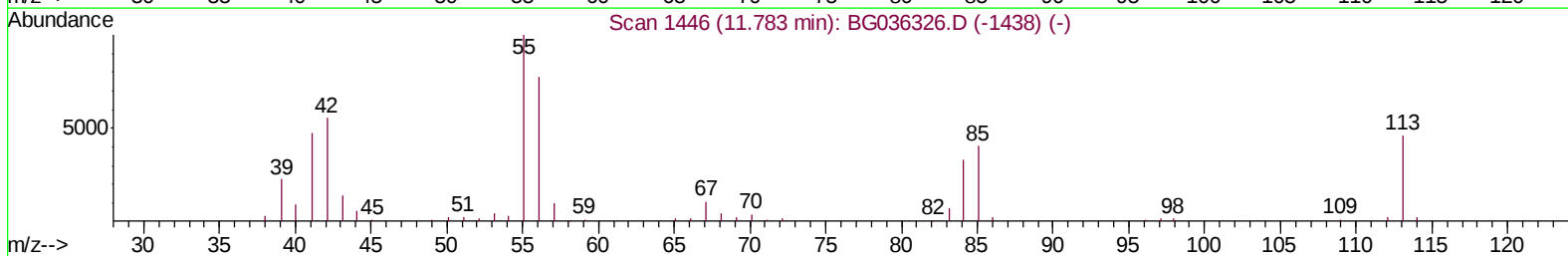
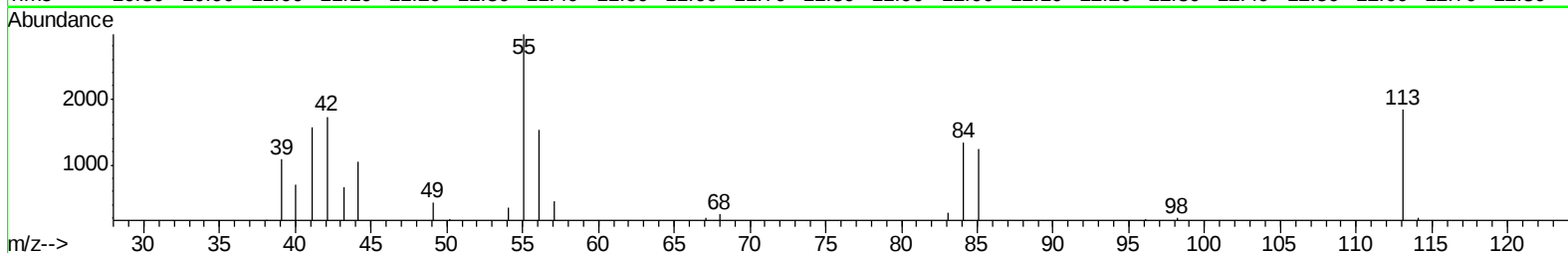
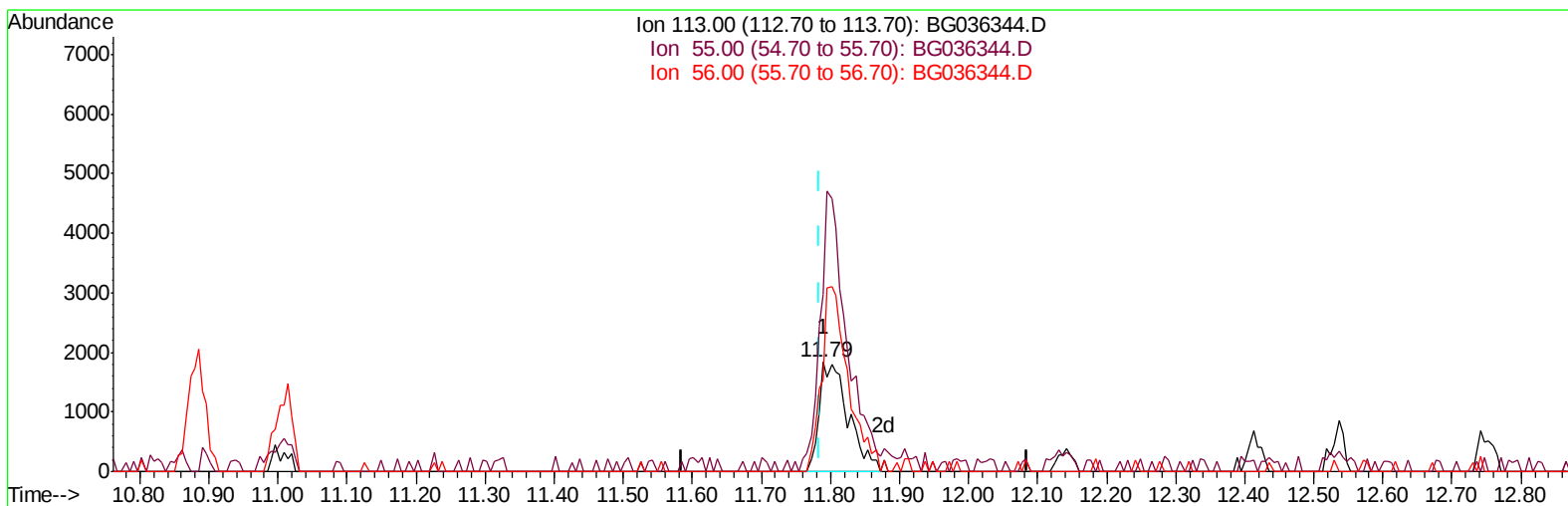
11.790min (+0.005) 1.25ng/ul

response 1803

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 162.10 |
| 56.00  | 115.90 | 83.31# |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

(32) Caprolactam

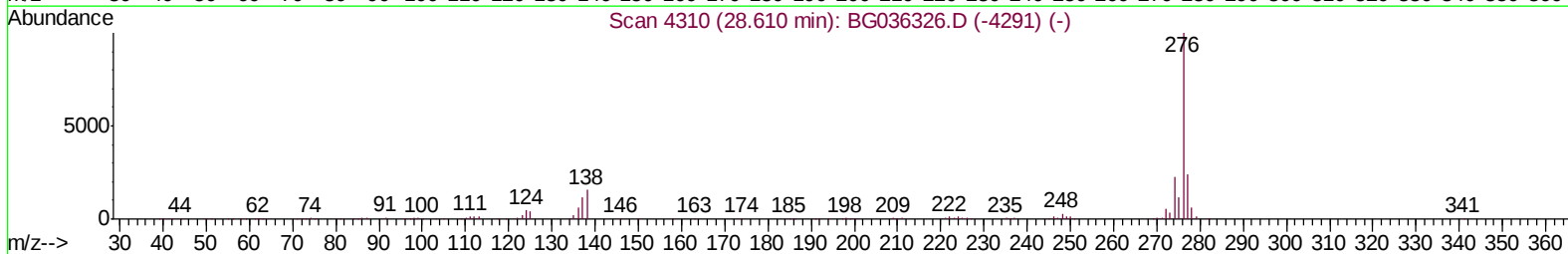
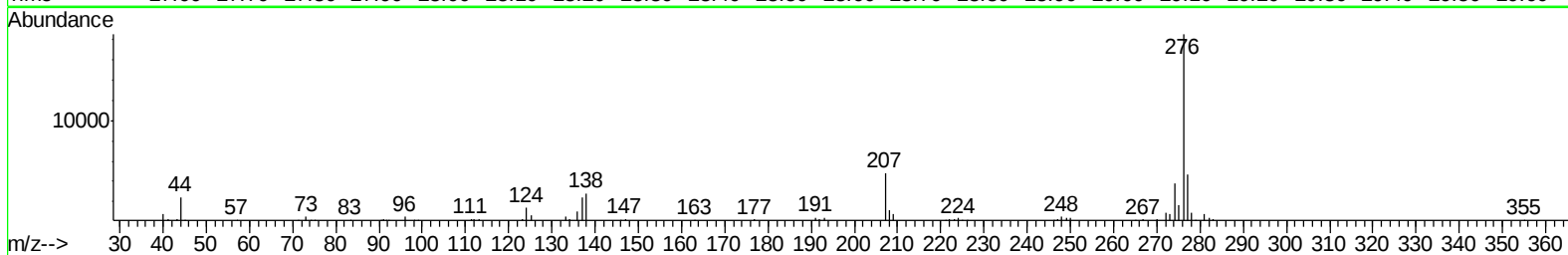
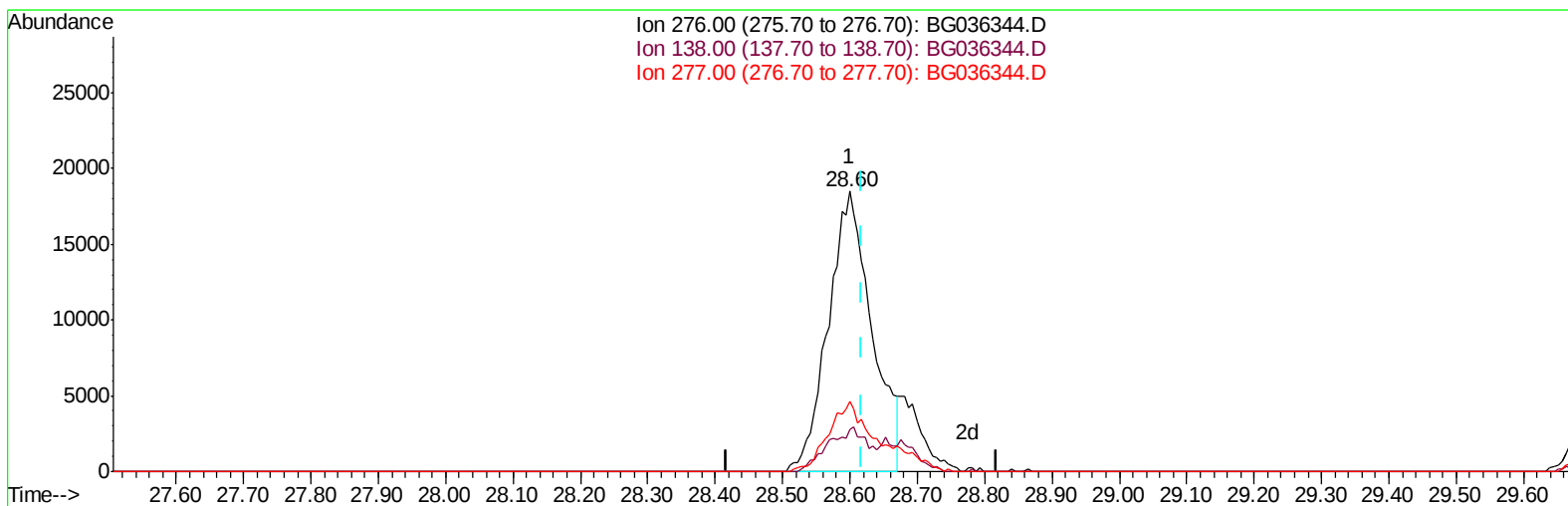
11.790min (+0.005) 3.69ng/ul m

response 5321

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 162.10 |
| 56.00  | 115.90 | 83.31# |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

(92) Indeno(1,2,3-cd)pyrene

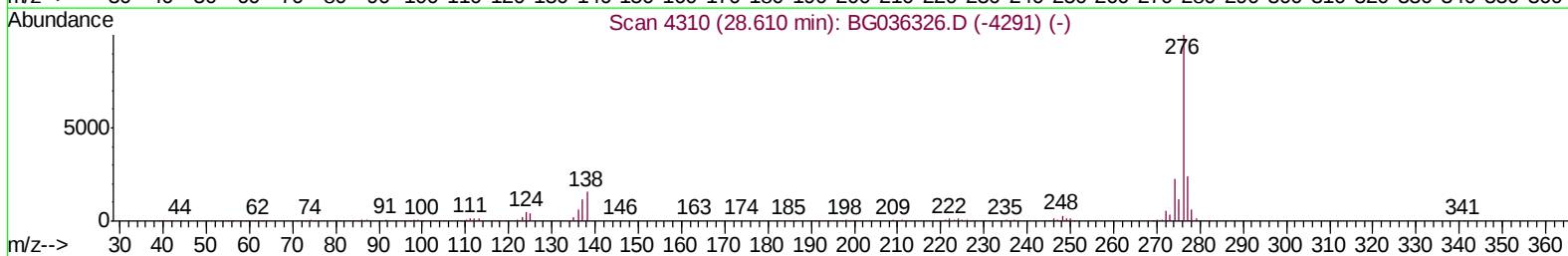
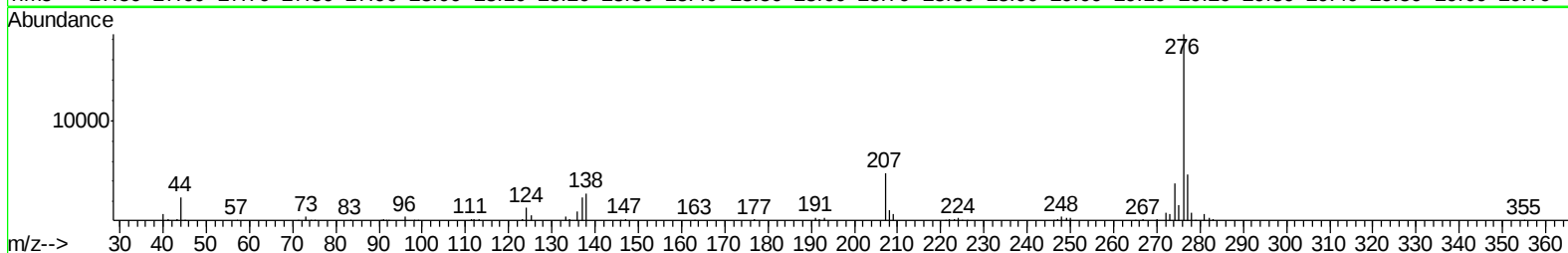
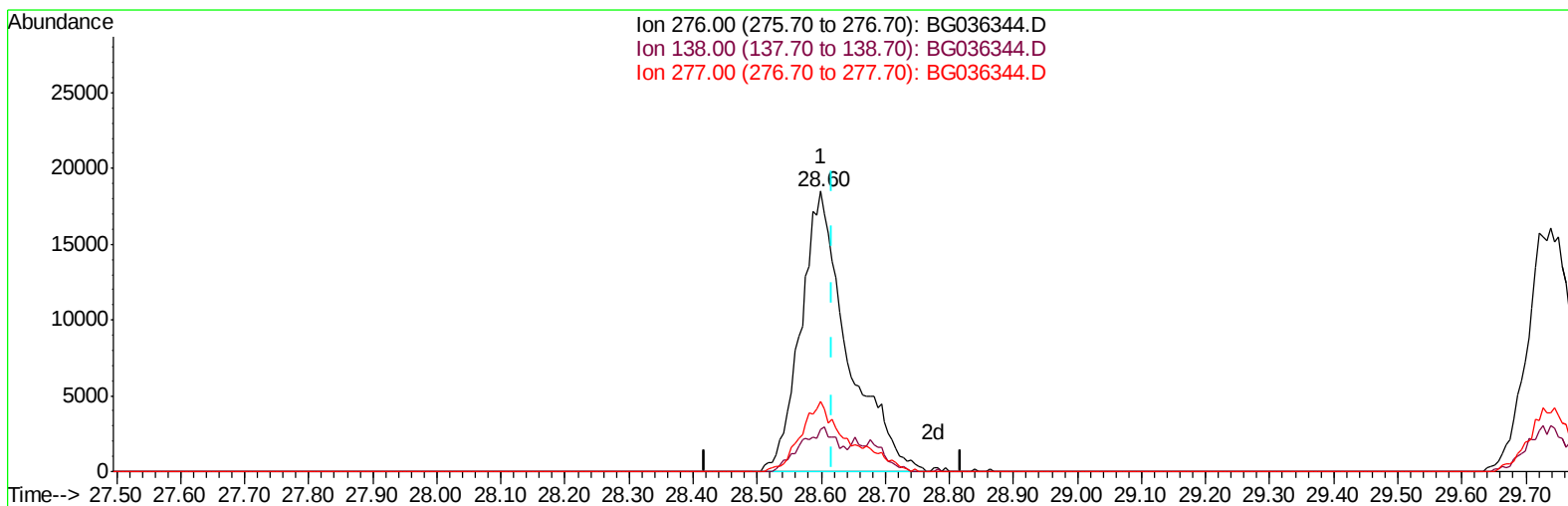
28.599min (-0.019) 3.31ng/ul

response 82997

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 13.60 | 15.12 |
| 277.00 | 23.60 | 25.04 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

(92) Indeno(1,2,3-cd)pyrene

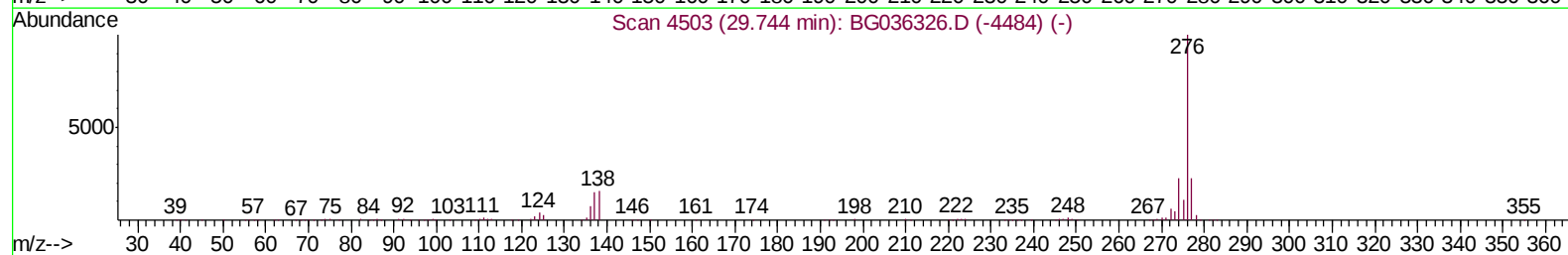
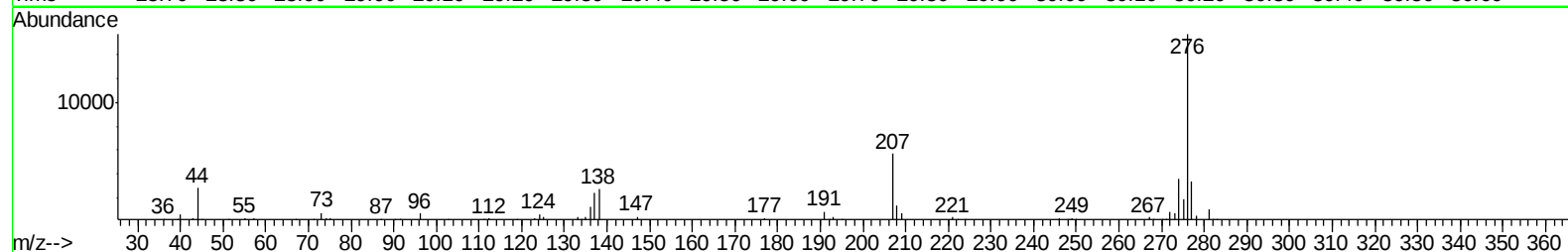
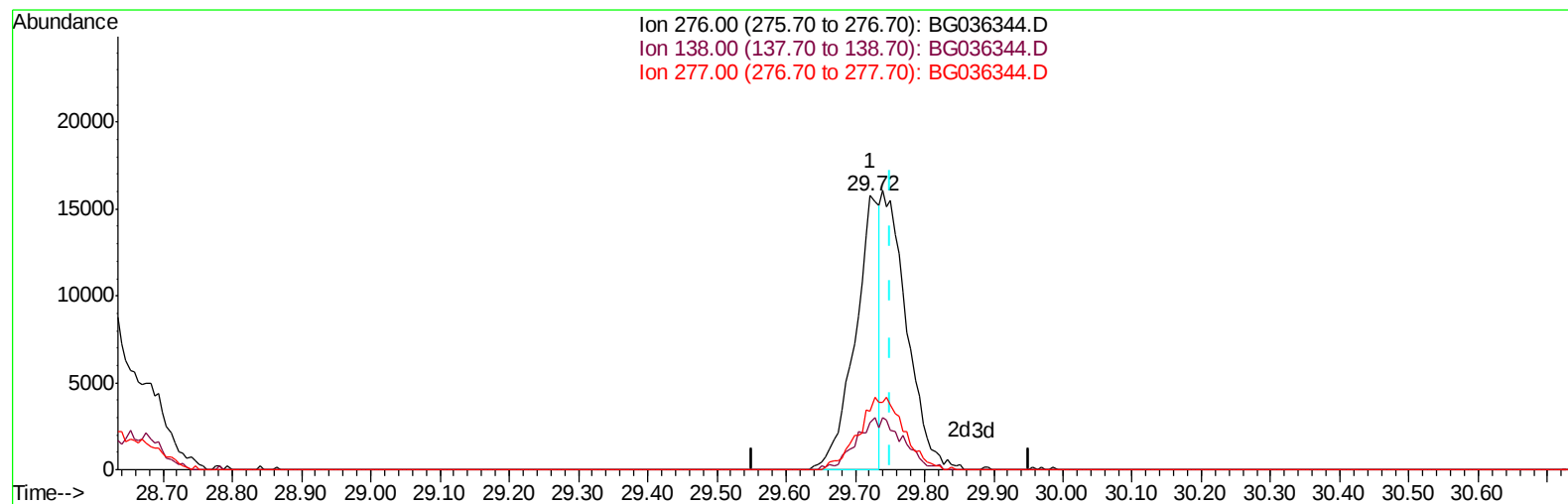
28.599min (-0.019) 3.77ng/ul m

response 94466

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 13.60 | 15.12 |
| 277.00 | 23.60 | 25.04 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

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

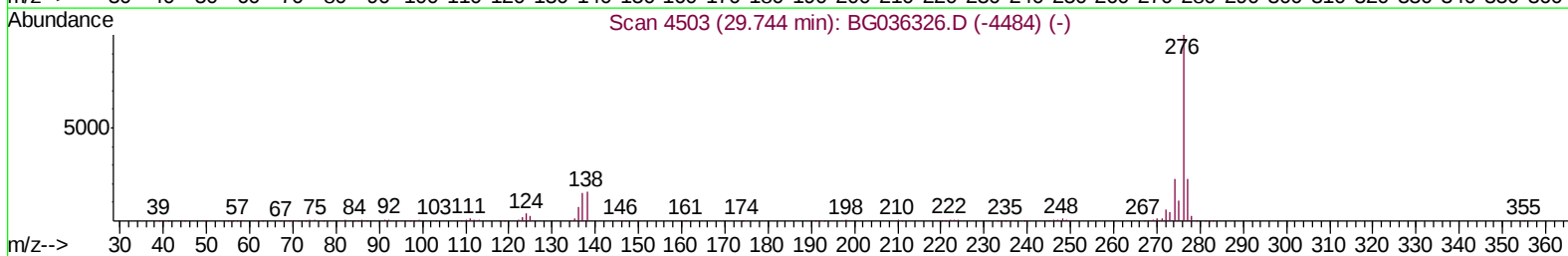
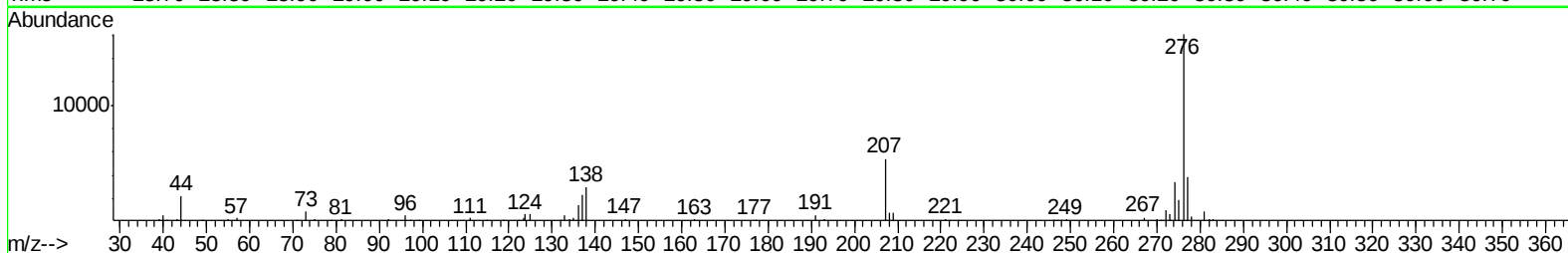
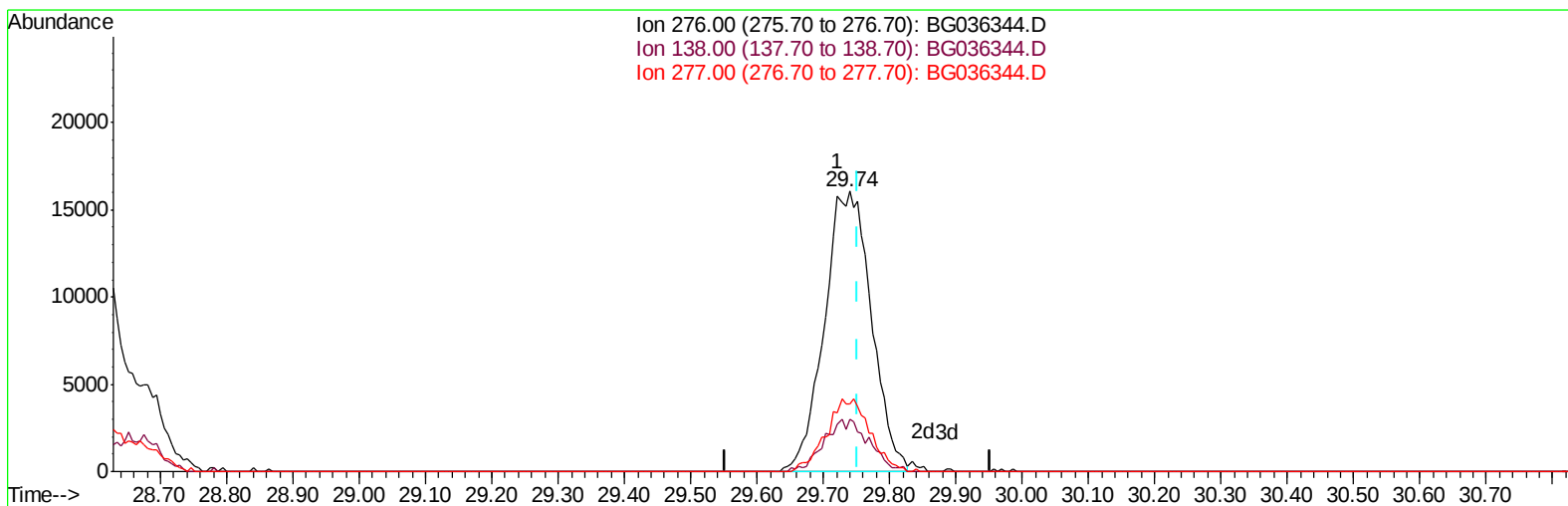
29.722min (-0.030) 1.84ng/ul

response 38067

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 276.00 | 100   | 100    |
| 138.00 | 14.00 | 17.01# |
| 277.00 | 23.80 | 21.48  |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
Data File : BG036344.D  
Acq On : 22 Aug 2018 15:58  
Operator : JU/SJ  
Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:32:48 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration



TIC: BG036344.D

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

29.739min (-0.013) 3.80ng/ul m

response 78511

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 276.00 | 100   | 100    |
| 138.00 | 14.00 | 18.81# |
| 277.00 | 23.80 | 24.14  |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036344.D  
 Acq On : 22 Aug 2018 15:58  
 Operator : JU/SJ  
 Sample : MDL-ME-04  
 Misc : 1PPM/4PPM  
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:36:24 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:52:18 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.08  | 152  | 50418    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.88 | 136  | 223603   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.70 | 164  | 140731   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.44 | 188  | 349073   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.71 | 240  | 375677   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.94 | 264  | 375931   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.51  | 96  | 8410   | 6.21  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.22  | 99  | 153815 | 32.79 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.40  | 67  | 105845 | 32.94 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.60  | 132 | 108817 | 32.13 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.76  | 113 | 122102 | 34.10 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.23  | 128 | 52562  | 38.53 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.96  | 143 | 53876  | 39.26 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.49 | 165 | 116814 | 31.80 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.01 | 131 | 157546 | 38.77 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.10 | 166 | 396519 | 33.70 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.39 | 160 | 478832 | 33.75 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.87 | 143 | 60760  | 34.91 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.69 | 176 | 335685 | 32.80 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.79 | 200 | 39786  | 31.47 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.54 | 188 | 540422 | 33.18 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.82 | 212 | 629992 | 32.73 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.72 | 264 | 610655 | 29.93 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |       |              | Qvalue |
|--------------------------------|-------|-----|-------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.55  | 88  | 2731  | 1.912 ng/uL# | 11     |
| 4) Benzaldehyde                | 7.21  | 77  | 6804  | 2.164 ng/ul# | 85     |
| 6) Phenol                      | 7.25  | 94  | 19128 | 4.080 ng/ul# | 90     |
| 8) Bis(2-Chloroethyl)ether     | 7.49  | 93  | 14832 | 4.260 ng/ul  | 87     |
| 10) 2-Chlorophenol             | 7.64  | 128 | 13134 | 3.877 ng/ul  | 98     |
| 11) 2-Methylphenol             | 8.51  | 108 | 13657 | 3.824 ng/ul  | 95     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.61  | 45  | 29904 | 3.902 ng/ul# | 84     |
| 14) Acetophenone               | 8.89  | 105 | 22805 | 4.155 ng/ul# | 85     |
| 15) N-Nitroso-di-n-propylamine | 8.88  | 70  | 12810 | 3.874 ng/ul# | 88     |
| 16) 4-Methylphenol             | 8.83  | 108 | 13729 | 3.674 ng/ul  | 99     |
| 17) Hexachloroethane           | 9.16  | 117 | 5333  | 3.951 ng/ul  | 89     |
| 20) Nitrobenzene               | 9.27  | 77  | 16395 | 4.236 ng/ul  | 91     |
| 21) Isophorone                 | 9.80  | 82  | 36451 | 3.880 ng/ul  | 95     |
| 23) 2-Nitrophenol              | 9.99  | 139 | 6666  | 4.589 ng/ul  | 93     |
| 24) 2,4-Dimethylphenol         | 10.04 | 107 | 16675 | 3.758 ng/ul  | 93     |
| 25) Bis(2-Chloroethoxy)methane | 10.29 | 93  | 19166 | 3.773 ng/ul  | 95     |
| 27) 2,4-Dichlorophenol         | 10.51 | 162 | 13494 | 3.877 ng/ul# | 80     |
| 28) Naphthalene                | 10.93 | 128 | 46868 | 4.090 ng/ul  | 98     |
| 30) 4-Chloroaniline            | 11.03 | 127 | 15575 | 3.934 ng/ul  | 93     |
| 31) Hexachlorobutadiene        | 11.22 | 225 | 10120 | 4.074 ng/ul  | 95     |
| 32) Caprolactam                | 11.79 | 113 | 5321m | 3.695 ng/ul  |        |
| 33) 4-Chloro-3-methylphenol    | 12.14 | 107 | 15459 | 3.781 ng/ul  | 98     |
| 34) 2-Methylnaphthalene        | 12.53 | 142 | 34368 | 4.018 ng/ul  | 100    |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
 Data File : BG036344.D  
 Acq On : 22 Aug 2018 15:58  
 Operator : JU/SJ  
 Sample : MDL-ME-04  
 Misc : 1PPM/4PPM  
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:36:24 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 10:52:18 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.75 | 142  | 34026    | 4.079 | ng/ul# | 90       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.89 | 216  | 18915    | 3.946 | ng/ul# | 96       |
| 38) Hexachlorocyclopentadiene  | 12.88 | 237  | 9951     | 3.469 | ng/ul# | 94       |
| 39) 2,4,6-Trichlorophenol      | 13.13 | 196  | 11225    | 3.791 | ng/ul  | 89       |
| 40) 2,4,5-Trichlorophenol      | 13.19 | 196  | 12107    | 3.945 | ng/ul  | 95       |
| 41) 1,1'-Biphenyl              | 13.53 | 154  | 44881    | 4.075 | ng/ul  | 95       |
| 42) 2-Chloronaphthalene        | 13.58 | 162  | 35189    | 4.164 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 13.77 | 65   | 8599     | 3.889 | ng/ul  | 93       |
| 45) Dimethylphthalate          | 14.15 | 163  | 45492    | 4.029 | ng/ul# | 98       |
| 46) 2,6-Dinitrotoluene         | 14.26 | 165  | 6383     | 4.045 | ng/ul# | 91       |
| 48) Acenaphthylene             | 14.42 | 152  | 53217    | 3.979 | ng/ul  | 99       |
| 49) 3-Nitroaniline             | 14.59 | 138  | 6270     | 3.813 | ng/ul# | 87       |
| 50) Acenaphthene               | 14.76 | 153  | 37652    | 3.932 | ng/ul  | 98       |
| 51) 2,4-Dinitrophenol          | 14.79 | 184  | 2374     | 2.958 | ng/ul# | 90       |
| 53) 4-Nitrophenol              | 14.88 | 109  | 6235     | 3.860 | ng/ul# | 84       |
| 54) Dibenzofuran               | 15.09 | 168  | 54363    | 4.102 | ng/ul  | 99       |
| 55) 2,4-Dinitrotoluene         | 15.04 | 165  | 8961     | 4.106 | ng/ul# | 77       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.32 | 232  | 10637    | 3.701 | ng/ul# | 95       |
| 57) Diethylphthalate           | 15.51 | 149  | 45867    | 4.003 | ng/ul  | 98       |
| 59) Fluorene                   | 15.74 | 166  | 43298    | 4.052 | ng/ul  | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.74 | 204  | 23415    | 4.022 | ng/ul  | 96       |
| 61) 4-Nitroaniline             | 15.74 | 138  | 6053     | 2.906 | ng/ul# | 81       |
| 64) 4,6-Dinitro-2-methylphenol | 15.81 | 198  | 4699     | 3.598 | ng/ul# | 66       |
| 65) N-Nitrosodiphenylamine     | 15.95 | 169  | 38130    | 3.899 | ng/ul  | 94       |
| 66) 4-Bromophenyl-phenylether  | 16.63 | 248  | 14771    | 3.826 | ng/ul  | 97       |
| 67) Hexachlorobenzene          | 16.74 | 284  | 14872    | 3.859 | ng/ul  | 92       |
| 68) Atrazine                   | 16.90 | 200  | 15460    | 3.833 | ng/ul  | 94       |
| 69) Pentachlorophenol          | 17.08 | 266  | 7597     | 3.251 | ng/ul  | 88       |
| 70) Phenanthrene               | 17.48 | 178  | 74111    | 4.085 | ng/ul  | 100      |
| 72) Anthracene                 | 17.58 | 178  | 74884    | 4.102 | ng/ul  | 95       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.50 | 216  | 20816    | 3.903 | ng/uL  | 95       |
| 74) Pentachlorobenzene         | 15.02 | 250  | 18387    | 3.960 | ng/uL  | 98       |
| 75) Carbazole                  | 17.84 | 167  | 65054    | 4.239 | ng/ul# | 94       |
| 76) Di-n-butylphthalate        | 18.41 | 149  | 75702    | 3.776 | ng/ul  | 100      |
| 77) Fluoranthene               | 19.49 | 202  | 91231    | 4.337 | ng/ul# | 95       |
| 80) Pyrene                     | 19.85 | 202  | 91473    | 4.018 | ng/ul# | 93       |
| 81) Butylbenzylphthalate       | 20.74 | 149  | 32308    | 3.487 | ng/ul  | 88       |
| 82) 3,3'-Dichlorobenzidine     | 21.60 | 252  | 23435    | 3.028 | ng/ul  | 95       |
| 83) Benzo(a)anthracene         | 21.69 | 228  | 95096    | 4.006 | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.63 | 149  | 47358    | 3.585 | ng/ul# | 97       |
| 85) Chrysene                   | 21.75 | 228  | 88366    | 4.058 | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.85 | 149  | 77572    | 3.671 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.91 | 252  | 88784    | 3.988 | ng/ul# | 96       |
| 89) Benzo(k)fluoranthene       | 23.98 | 252  | 85449    | 3.981 | ng/ul# | 98       |
| 91) Benzo(a)pyrene             | 24.79 | 252  | 81290    | 3.859 | ng/ul# | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 28.60 | 276  | 94466m   | 3.766 | ng/ul  |          |
| 93) Dibenzo(a,h)anthracene     | 28.67 | 278  | 78510    | 3.860 | ng/ul# | 94       |
| 94) Benzo(g,h,i)perylene       | 29.74 | 276  | 78511m   | 3.796 | ng/ul  |          |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
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Sample : MDL-ME-04  
Misc : 1PPM/4PPM  
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 22 16:36:24 2018  
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Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 22 10:52:18 2018  
Response via : Initial Calibration

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082218\  
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