

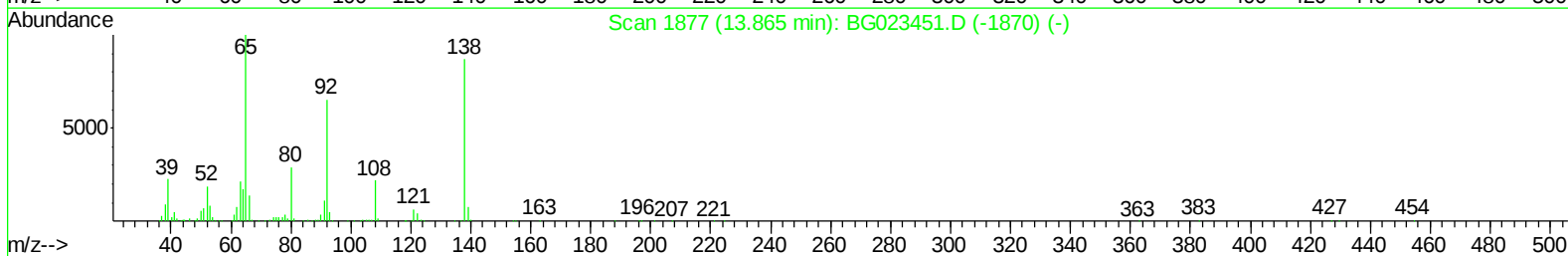
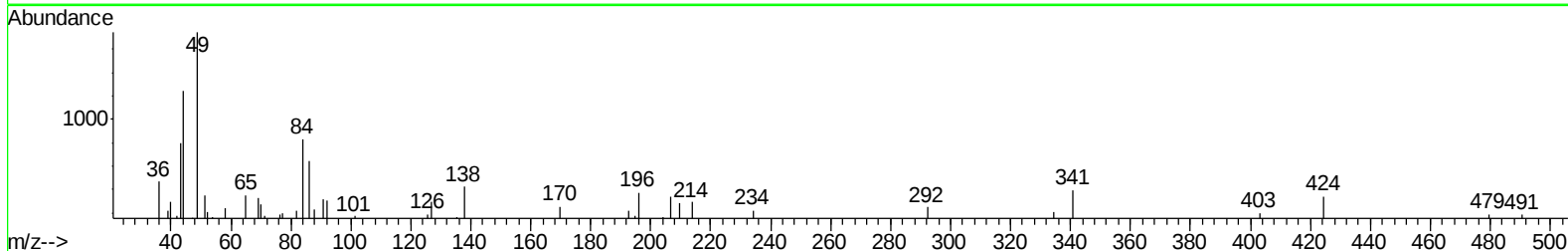
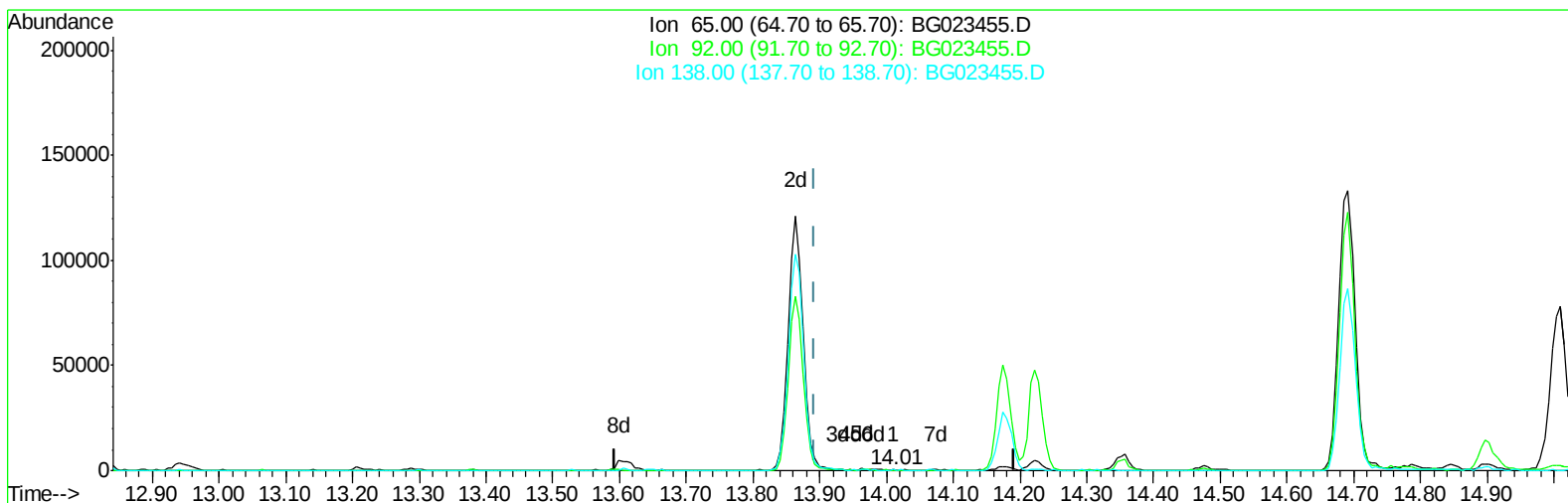
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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8/5/2016 6:59:57 PM

Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(42) 2-Nitroaniline

14.010min (+0.118) 0.03ng/ul

response 282

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 65.00  | 100    | 100    |
| 92.00  | 83.90  | 86.12  |
| 138.00 | 137.00 | 119.26 |
| 0.00   | 0.00   | 0.00   |

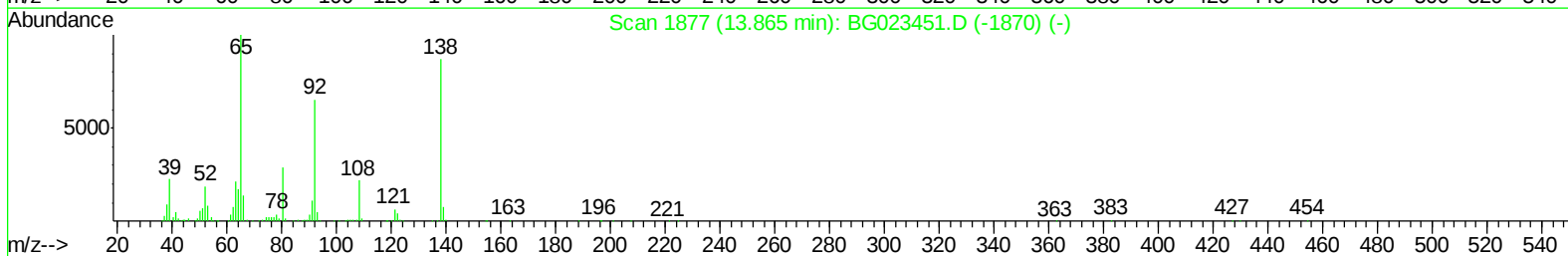
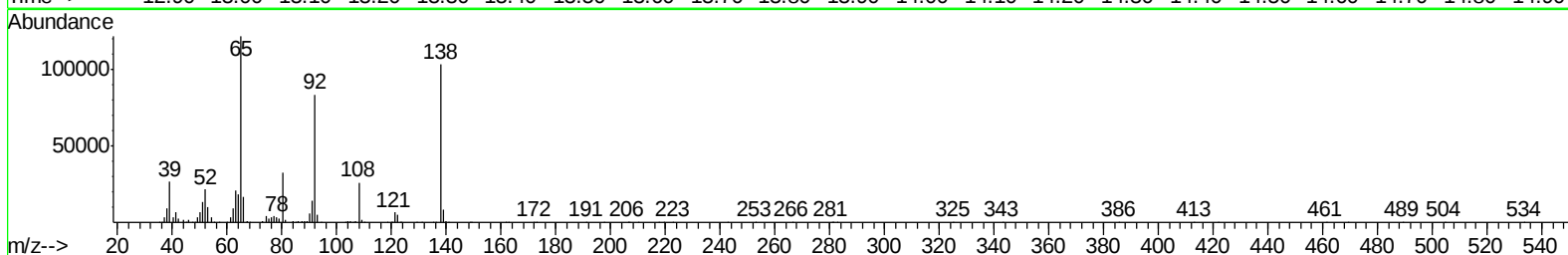
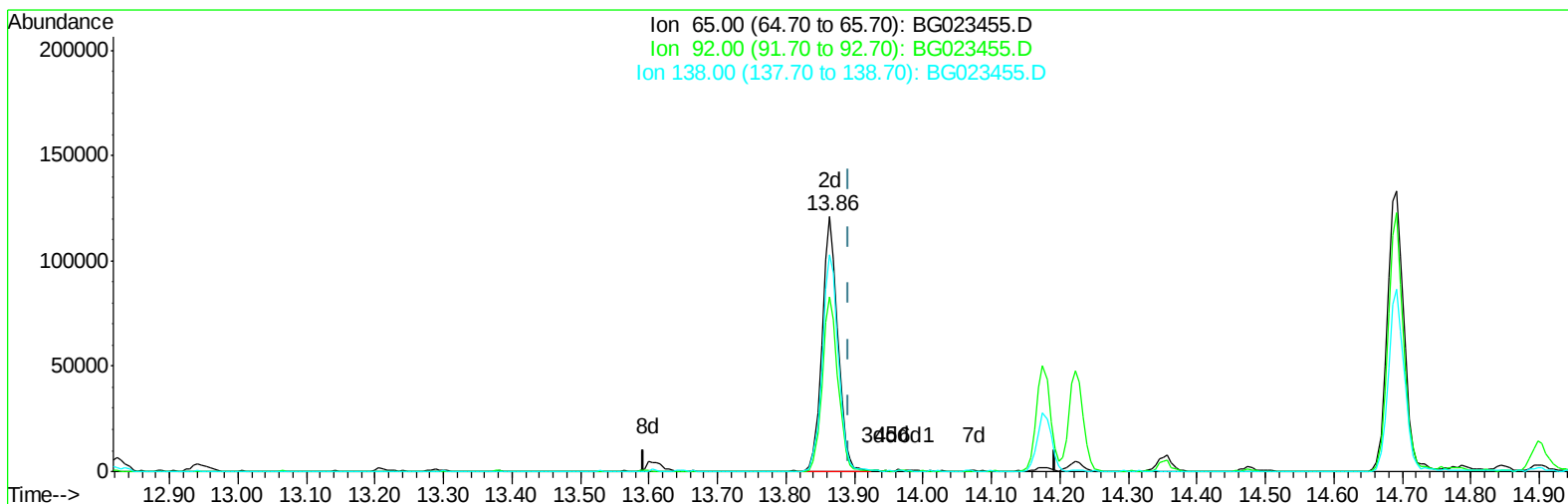
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Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
APPROVED

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8/5/2016 6:59:57 PM

Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(42) 2-Nitroaniline

13.863min (-0.029) 22.18ng/ul m

response 193914

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 65.00  | 100    | 100    |
| 92.00  | 83.90  | 68.55  |
| 138.00 | 137.00 | 84.90# |
| 0.00   | 0.00   | 0.00   |

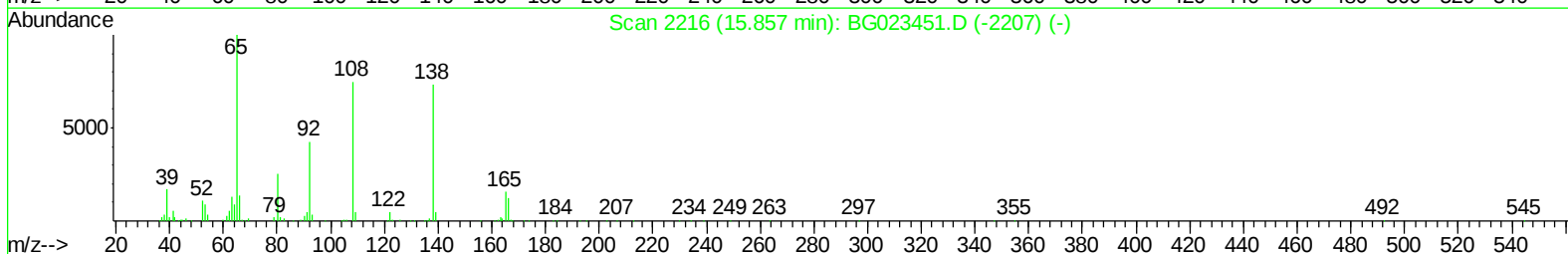
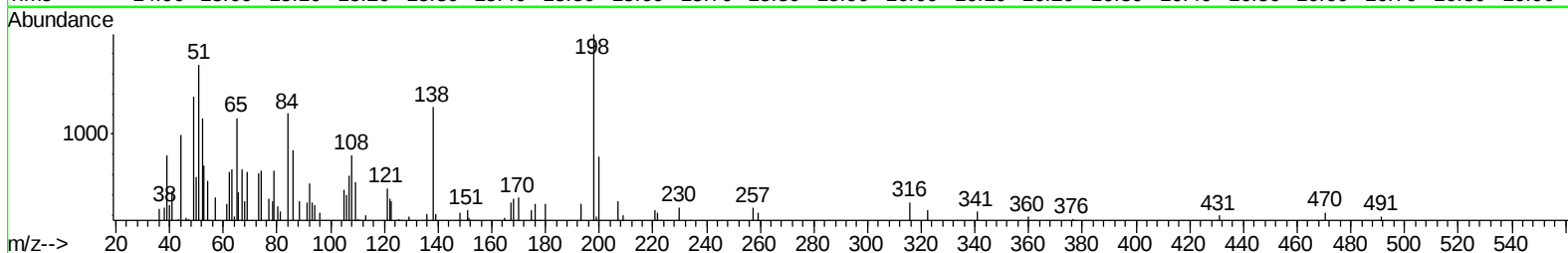
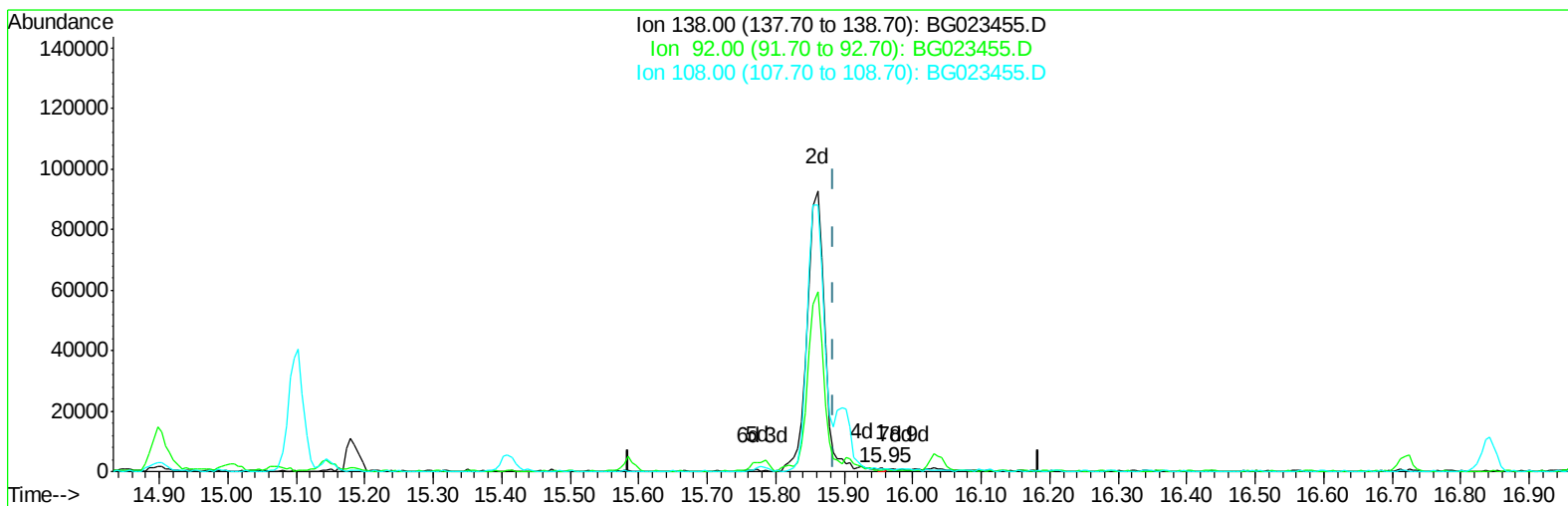
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(60) 4-Nitroaniline

15.955min (+0.070) 0.07ng/ul

response 531

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 48.70 | 40.88 |
| 108.00 | 61.50 | 62.59 |
| 0.00   | 0.00  | 0.00  |

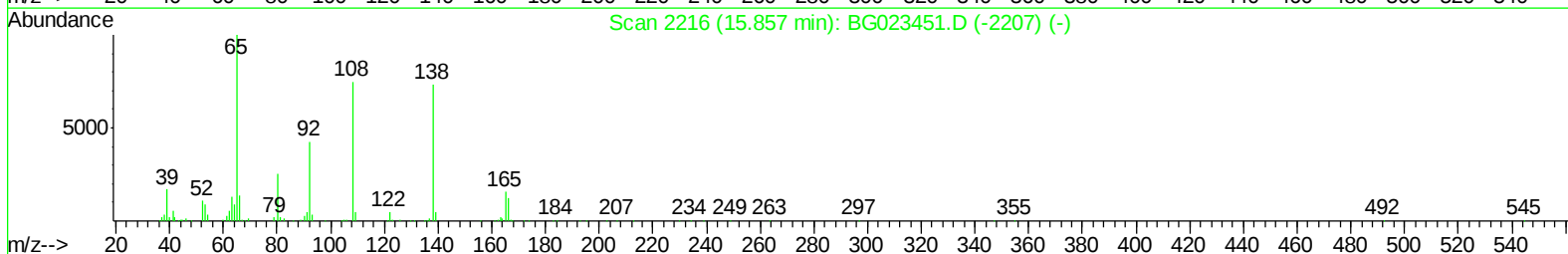
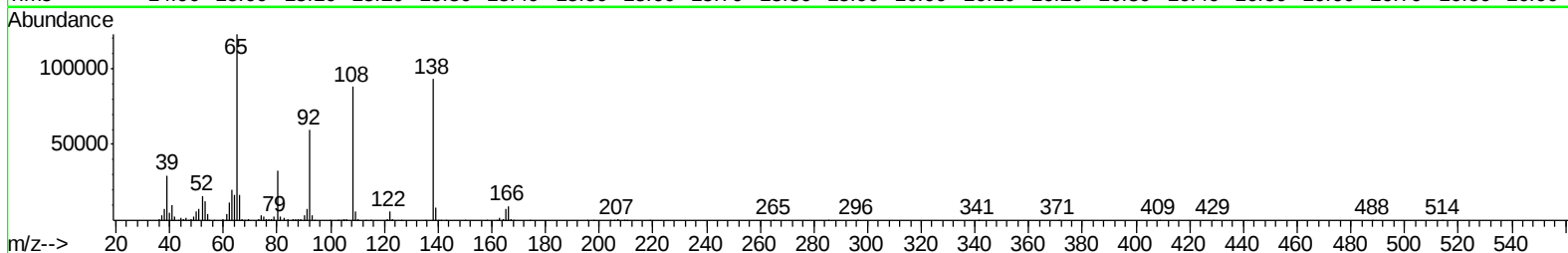
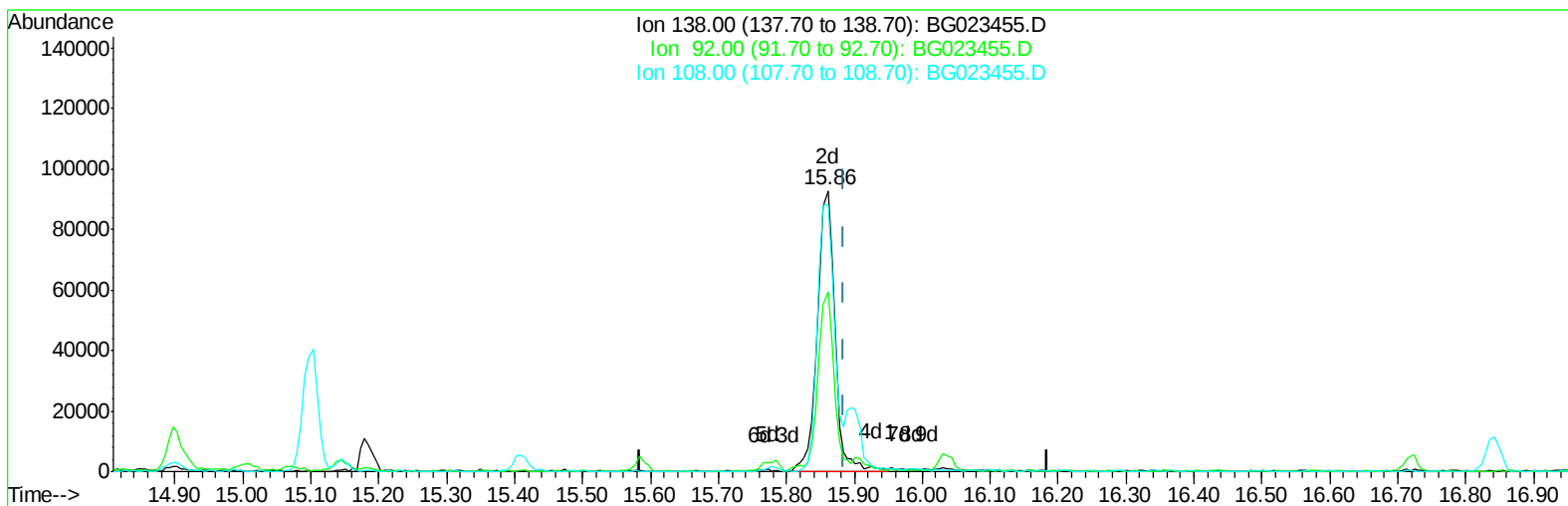
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(60) 4-Nitroaniline

15.861min (-0.024) 22.94ng/ul m

response 164799

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 138.00 | 100   | 100    |
| 92.00  | 48.70 | 64.16# |
| 108.00 | 61.50 | 94.95# |
| 0.00   | 0.00  | 0.00   |

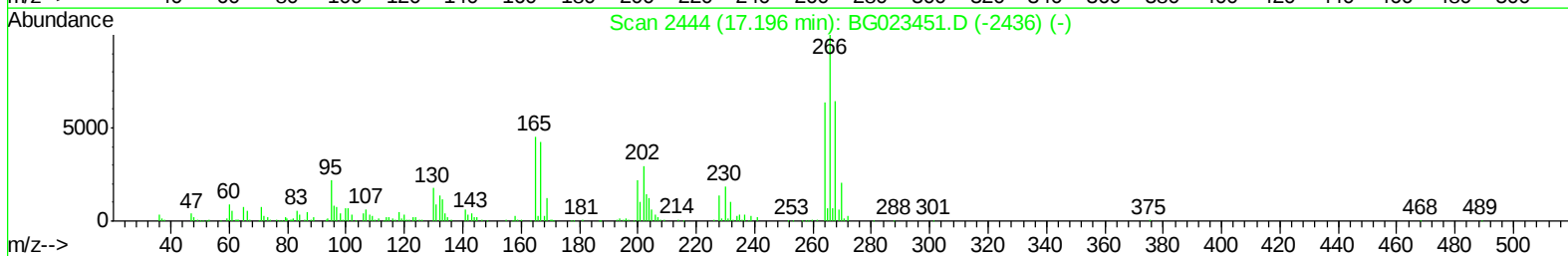
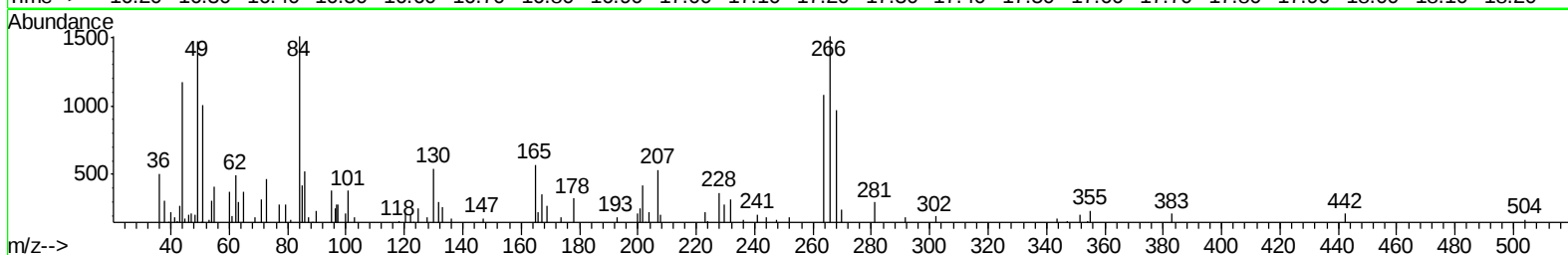
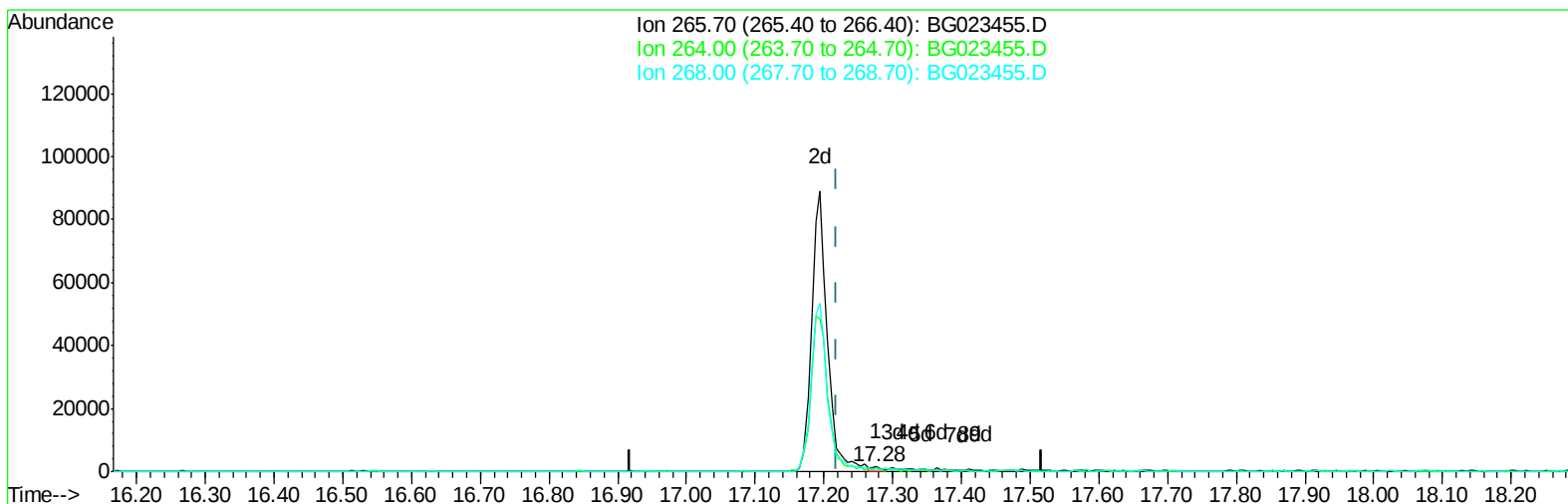
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(68) Pentachlorophenol (C)

17.276min (+0.058) 0.20ng/ul

response 1356

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 265.70 | 100   | 100   |
| 264.00 | 68.40 | 71.58 |
| 268.00 | 72.30 | 64.28 |
| 0.00   | 0.00  | 0.00  |

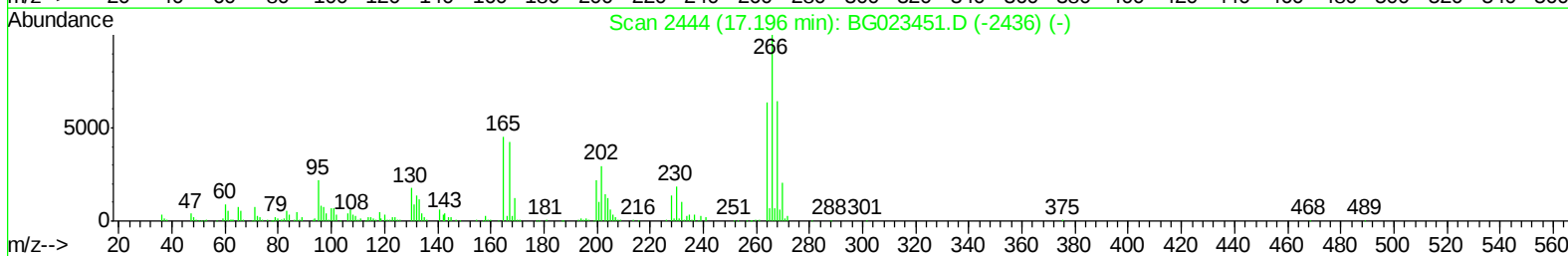
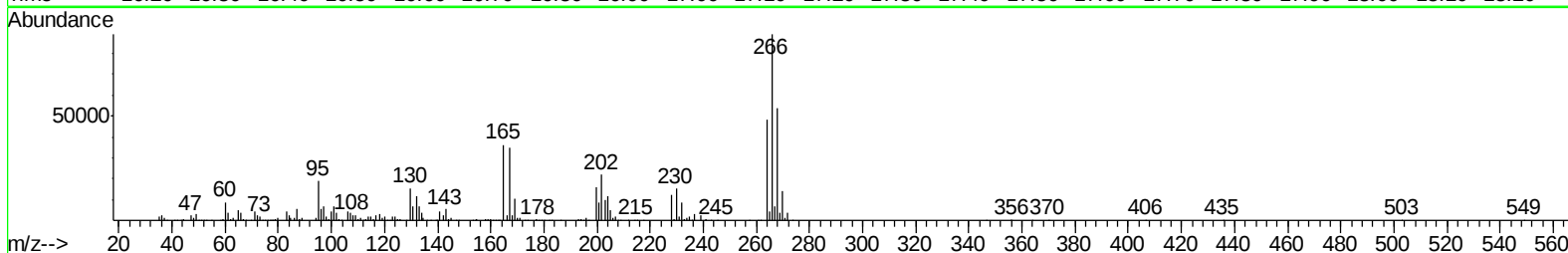
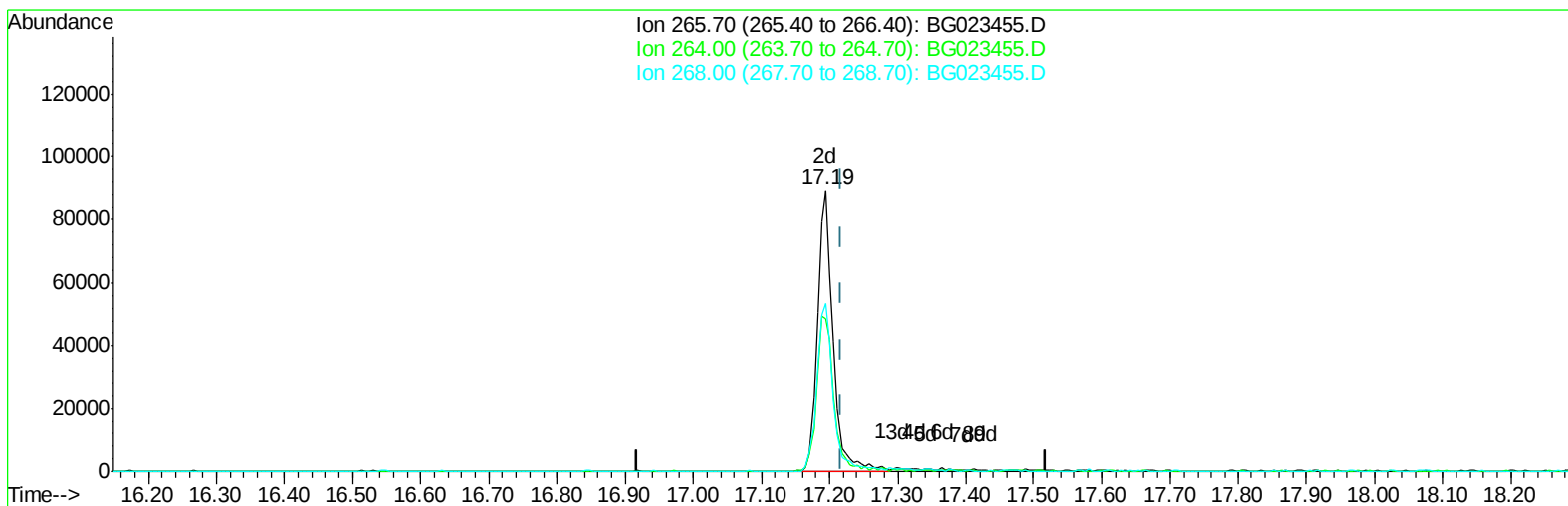
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(68) Pentachlorophenol (C)

17.194min (-0.024) 21.39ng/ul m

response 143969

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 265.70 | 100   | 100    |
| 264.00 | 68.40 | 54.40# |
| 268.00 | 72.30 | 60.08  |
| 0.00   | 0.00  | 0.00   |

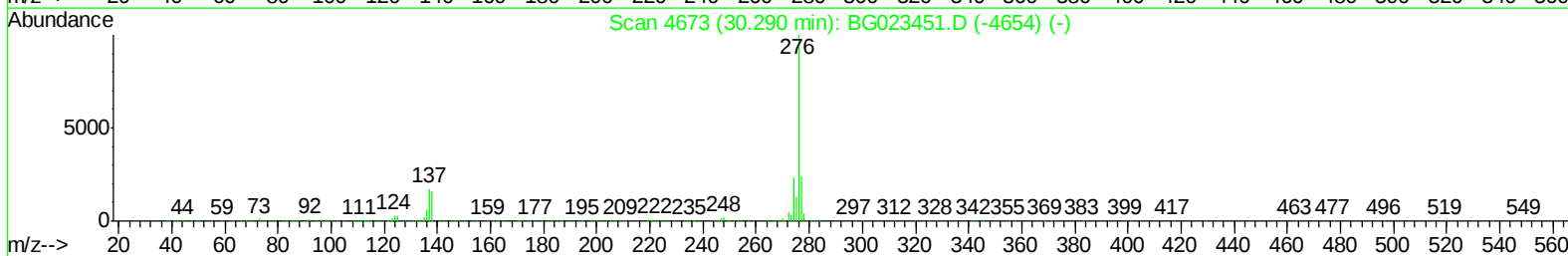
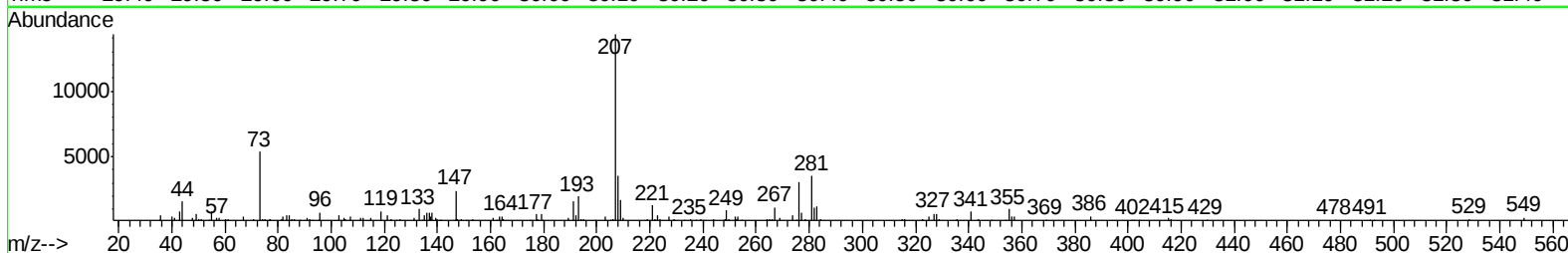
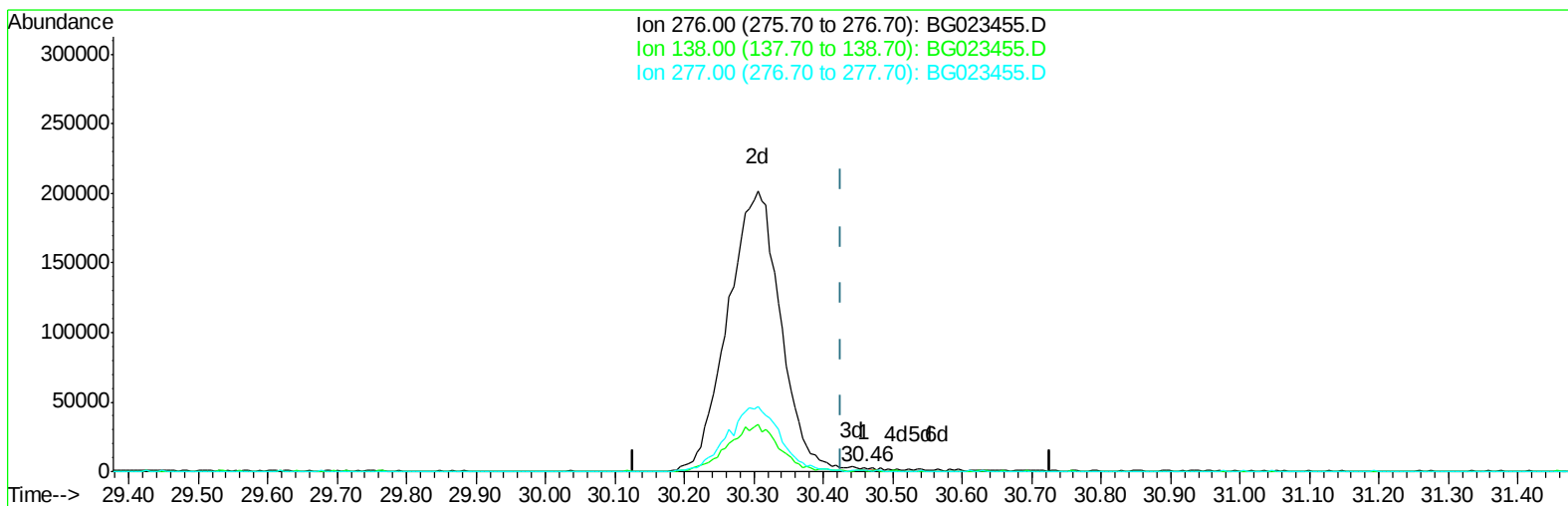
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Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

(91) Benzo(g,h,i)perylene  
30.459min (+0.032) 0.04ng/ul  
response 2126

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 27.60 | 23.74 |
| 277.00 | 21.20 | 25.08 |
| 0.00   | 0.00  | 0.00  |

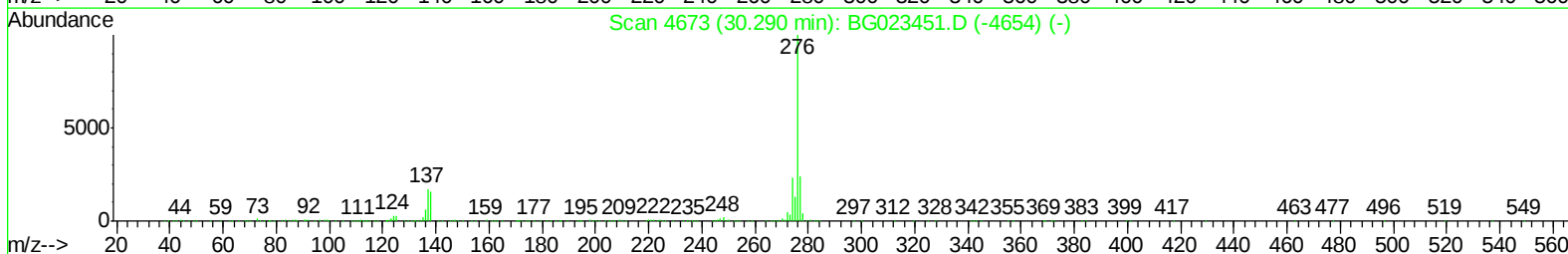
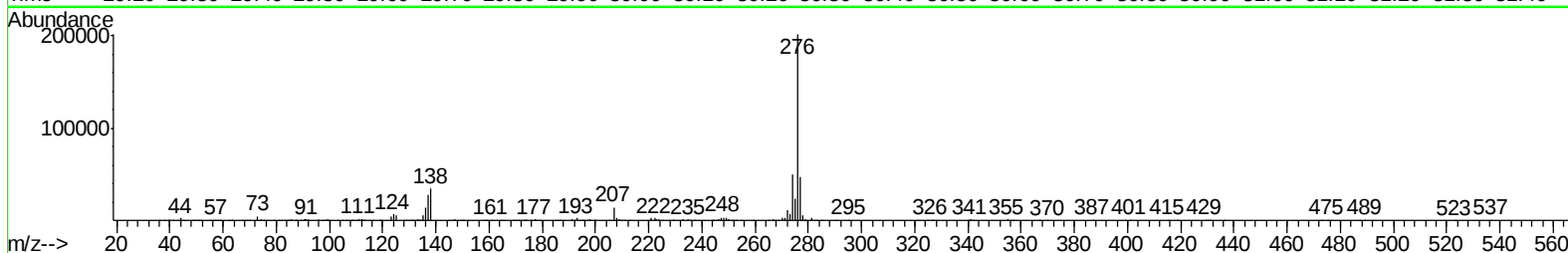
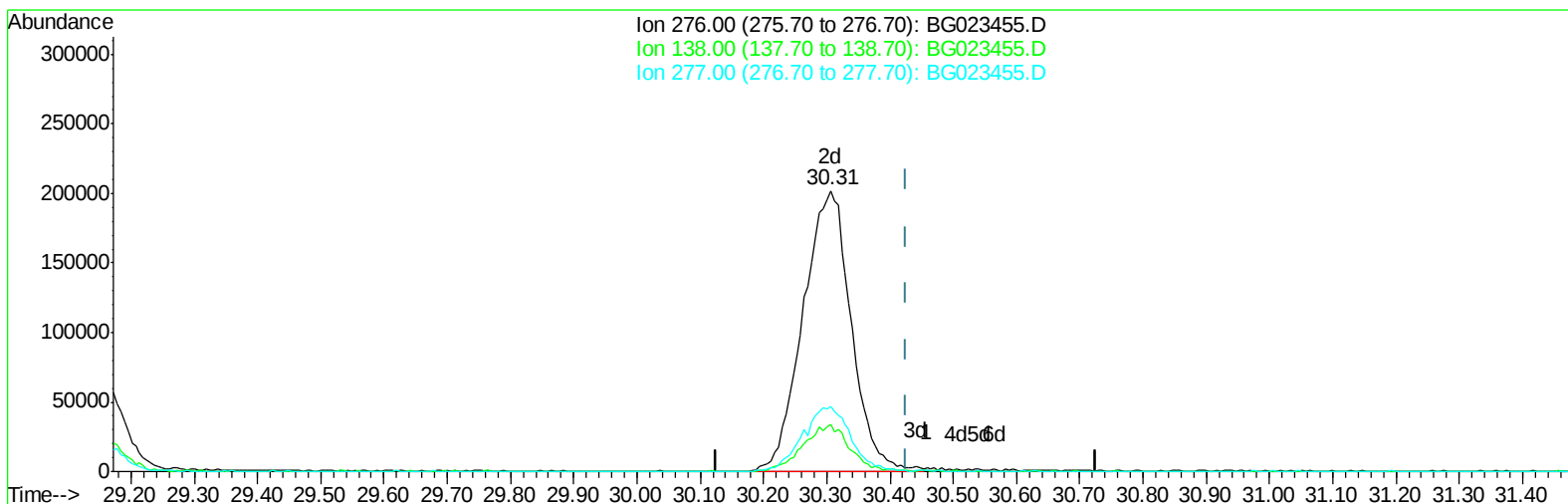
Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
Data File : BG023455.D  
Acq On : 4 Aug 2016 17:08  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02007

Manual Integrations  
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Quant Time: Aug 04 17:41:22 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 04 17:27:34 2016  
Response via : Initial Calibration



TIC: BG023455.D

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

30.306min (-0.120) 21.33ng/ul m

response 1067881

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 276.00 | 100   | 100    |
| 138.00 | 27.60 | 16.92# |
| 277.00 | 21.20 | 23.22  |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
 Data File : BG023455.D  
 Acq On : 4 Aug 2016 17:08  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02007

Manual Integrations  
 APPROVED

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 8/5/2016 6:59:57 PM

Quant Time: Aug 04 17:46:07 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 04 17:27:34 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.12  | 152  | 106579   | 20.00 | ng/ul | -0.02    |
| 18) Naphthalene-d8        | 10.95 | 136  | 494095   | 20.00 | ng/ul | -0.03    |
| 35) Acenaphthene-d10      | 14.79 | 164  | 374765   | 20.00 | ng/ul | -0.02    |
| 61) Phenanthrene-d10      | 17.54 | 188  | 998307   | 20.00 | ng/ul | -0.03    |
| 75) Chrysene-d12          | 21.86 | 240  | 989491   | 20.00 | ng/ul | -0.04    |
| 83) Perylene-d12          | 25.24 | 264  | 935889   | 20.00 | ng/ul | -0.07    |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.49  | 96  | 19971   | 7.97  | ng/uL | -0.02 |
| 5) Phenol-d5                   | 7.27  | 99  | 224277  | 20.89 | ng/ul | -0.02 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.43  | 67  | 147368  | 21.42 | ng/ul | -0.02 |
| 9) 2-Chlorophenol-d4           | 7.65  | 132 | 153148  | 20.91 | ng/ul | -0.02 |
| 13) 4-Methylphenol-d8          | 8.83  | 113 | 175547  | 20.94 | ng/ul | -0.02 |
| 19) Nitrobenzene-d5            | 9.29  | 128 | 78104   | 20.05 | ng/ul | -0.03 |
| 22) 2-Nitrophenol-d4           | 10.02 | 143 | 97024   | 21.69 | ng/ul | -0.03 |
| 26) 2,4-Dichlorophenol-d3      | 10.57 | 165 | 187073  | 22.17 | ng/ul | -0.03 |
| 29) 4-Chloroaniline-d4         | 11.08 | 131 | 231662  | 23.40 | ng/ul | -0.03 |
| 43) Dimethylphthalate-d6       | 14.17 | 166 | 658025  | 21.44 | ng/ul | -0.03 |
| 46) Acenaphthylene-d8          | 14.47 | 160 | 744029  | 21.54 | ng/ul | -0.03 |
| 51) 4-Nitrophenol-d4           | 14.99 | 143 | 113639  | 21.69 | ng/ul | -0.02 |
| 57) Fluorene-d10               | 15.78 | 176 | 600918  | 21.20 | ng/ul | -0.03 |
| 62) 4,6-Dinitro-2-methylphenol | 15.90 | 200 | 135851  | 21.21 | ng/ul | -0.02 |
| 70) Anthracene-d10             | 17.64 | 188 | 956681  | 21.27 | ng/ul | -0.03 |
| 76) Pyrene-d10                 | 19.93 | 212 | 1079932 | 21.56 | ng/ul | -0.03 |
| 87) Benzo(a)pyrene-d12         | 25.00 | 264 | 915308  | 21.61 | ng/ul | -0.06 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.52  | 88  | 23885  | 9.08  | ng/uL# | 77     |
| 4) Benzaldehyde                | 7.25  | 77  | 152524 | 23.31 | ng/ul# | 79     |
| 6) Phenol                      | 7.30  | 94  | 236912 | 21.19 | ng/ul# | 73     |
| 8) Bis(2-Chloroethyl)ether     | 7.53  | 93  | 170533 | 21.01 | ng/ul  | 89     |
| 10) 2-Chlorophenol             | 7.68  | 128 | 153930 | 20.63 | ng/ul# | 84     |
| 11) 2-Methylphenol             | 8.56  | 108 | 174074 | 20.67 | ng/ul  | 96     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.65  | 45  | 241315 | 21.73 | ng/ul  | 97     |
| 14) Acetophenone               | 8.95  | 105 | 294556 | 21.91 | ng/ul# | 80     |
| 15) N-Nitroso-di-n-propylamine | 8.92  | 70  | 172621 | 21.34 | ng/ul# | 89     |
| 16) 4-Methylphenol             | 8.90  | 108 | 197046 | 20.88 | ng/ul  | 84     |
| 17) Hexachloroethane           | 9.21  | 117 | 63067  | 19.59 | ng/ul# | 75     |
| 20) Nitrobenzene               | 9.33  | 77  | 243924 | 21.37 | ng/ul# | 87     |
| 21) Isophorone                 | 9.86  | 82  | 460642 | 21.94 | ng/ul# | 93     |
| 23) 2-Nitrophenol              | 10.05 | 139 | 101998 | 21.12 | ng/ul# | 67     |
| 24) 2,4-Dimethylphenol         | 10.11 | 107 | 229755 | 21.68 | ng/ul  | 90     |
| 25) Bis(2-Chloroethoxy)methane | 10.34 | 93  | 256750 | 21.20 | ng/ul# | 97     |
| 27) 2,4-Dichlorophenol         | 10.60 | 162 | 184346 | 21.67 | ng/ul  | 98     |
| 28) Naphthalene                | 11.00 | 128 | 532193 | 21.22 | ng/ul  | 99     |
| 30) 4-Chloroaniline            | 11.11 | 127 | 229253 | 23.34 | ng/ul  | 97     |
| 31) Hexachlorobutadiene        | 11.28 | 225 | 121319 | 21.41 | ng/ul  | 87     |
| 32) Caprolactam                | 11.87 | 113 | 75486  | 21.36 | ng/ul# | 50     |
| 33) 4-Chloro-3-methylphenol    | 12.24 | 107 | 239787 | 21.90 | ng/ul# | 83     |
| 34) 2-Methylnaphthalene        | 12.61 | 142 | 432483 | 21.26 | ng/ul  | 96     |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080416\  
 Data File : BG023455.D  
 Acq On : 4 Aug 2016 17:08  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02007

Manual Integrations  
 APPROVED

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 8/5/2016 6:59:57 PM

Quant Time: Aug 04 17:46:07 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 04 17:27:34 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.97 | 216  | 244574   | 21.26 | ng/ul  | 97       |
| 37) Hexachlorocyclopentadiene  | 12.95 | 237  | 134007   | 21.06 | ng/ul  | 96       |
| 38) 2,4,6-Trichlorophenol      | 13.21 | 196  | 167734   | 21.05 | ng/ul  | 96       |
| 39) 2,4,5-Trichlorophenol      | 13.29 | 196  | 184095   | 21.55 | ng/ul  | 94       |
| 40) 1,1'-Biphenyl              | 13.61 | 154  | 589681   | 21.41 | ng/ul  | 99       |
| 41) 2-Chloronaphthalene        | 13.66 | 162  | 460962   | 21.52 | ng/ul  | 95       |
| 42) 2-Nitroaniline             | 13.86 | 65   | 193914m  | 22.18 | ng/ul  |          |
| 44) Dimethylphthalate          | 14.22 | 163  | 660876   | 21.75 | ng/ul  | 100      |
| 45) 2,6-Dinitrotoluene         | 14.36 | 165  | 138516   | 21.03 | ng/ul# | 83       |
| 47) Acenaphthylene             | 14.50 | 152  | 779403   | 21.58 | ng/ul  | 98       |
| 48) 3-Nitroaniline             | 14.69 | 138  | 137511   | 22.20 | ng/ul# | 59       |
| 49) Acenaphthene               | 14.84 | 153  | 518822   | 21.08 | ng/ul  | 99       |
| 50) 2,4-Dinitrophenol          | 14.90 | 184  | 86755    | 21.34 | ng/ul# | 88       |
| 52) 4-Nitrophenol              | 15.01 | 109  | 114189   | 21.76 | ng/ul# | 73       |
| 53) Dibenzofuran               | 15.18 | 168  | 786465   | 21.57 | ng/ul  | 88       |
| 54) 2,4-Dinitrotoluene         | 15.14 | 165  | 218564   | 21.90 | ng/ul# | 78       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.41 | 232  | 189869   | 21.63 | ng/ul  | 97       |
| 56) Diethylphthalate           | 15.58 | 149  | 690522   | 21.74 | ng/ul  | 97       |
| 58) Fluorene                   | 15.83 | 166  | 649365   | 21.38 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.82 | 204  | 326991   | 21.40 | ng/ul  | 99       |
| 60) 4-Nitroaniline             | 15.86 | 138  | 164799m  | 22.94 | ng/ul  |          |
| 63) 4,6-Dinitro-2-methylphenol | 15.91 | 198  | 142045   | 21.13 | ng/ul# | 85       |
| 64) N-Nitrosodiphenylamine     | 16.04 | 169  | 591094   | 21.19 | ng/ul  | 97       |
| 65) 4-Bromophenyl-phenylether  | 16.72 | 248  | 236937   | 20.70 | ng/ul  | 99       |
| 66) Hexachlorobenzene          | 16.84 | 284  | 258785   | 21.22 | ng/ul  | 94       |
| 67) Atrazine                   | 16.99 | 200  | 250434   | 21.73 | ng/ul  | 95       |
| 68) Pentachlorophenol          | 17.19 | 266  | 143969m  | 21.39 | ng/ul  |          |
| 69) Phenanthrene               | 17.59 | 178  | 1107483  | 21.42 | ng/ul  | 97       |
| 71) Anthracene                 | 17.68 | 178  | 1145664  | 21.72 | ng/ul  | 98       |
| 72) Carbazole                  | 17.95 | 167  | 1003336  | 23.60 | ng/ul  | 99       |
| 73) Di-n-butylphthalate        | 18.49 | 149  | 1177555  | 22.12 | ng/ul# | 95       |
| 74) Fluoranthene               | 19.60 | 202  | 1325226  | 25.04 | ng/ul# | 94       |
| 77) Pyrene                     | 19.96 | 202  | 1324078  | 21.65 | ng/ul  | 95       |
| 78) Butylbenzylphthalate       | 20.84 | 149  | 513363   | 21.88 | ng/ul# | 85       |
| 79) 3,3'-Dichlorobenzidine     | 21.75 | 252  | 390562   | 22.40 | ng/ul# | 97       |
| 80) Benzo(a)anthracene         | 21.83 | 228  | 1198219  | 21.50 | ng/ul  | 97       |
| 81) Bis(2-ethylhexyl)phthalate | 21.72 | 149  | 679597   | 21.78 | ng/ul# | 97       |
| 82) Chrysene                   | 21.91 | 228  | 1083318  | 21.38 | ng/ul  | 97       |
| 84) Di-n-octyl phthalate       | 22.98 | 149  | 1124688  | 23.79 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 24.16 | 252  | 1176524  | 22.10 | ng/ul# | 95       |
| 86) Benzo(k)fluoranthene       | 24.23 | 252  | 1127768  | 21.52 | ng/ul# | 92       |
| 88) Benzo(a)pyrene             | 25.07 | 252  | 1105182  | 21.50 | ng/ul# | 92       |
| 89) Indeno(1,2,3-cd)pyrene     | 29.09 | 276  | 1276990  | 21.15 | ng/ul# | 88       |
| 90) Dibenzo(a,h)anthracene     | 29.15 | 278  | 1109285  | 21.54 | ng/ul# | 93       |
| 91) Benzo(g,h,i)perylene       | 30.31 | 276  | 1067881m | 21.33 | ng/ul  |          |

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

