

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080718\  
 Data File : BG036176.D  
 Acq On : 8 Aug 2018 00:18  
 Operator : JU/SJ  
 Sample : J4303-07  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-A03

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG071218.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.127       | 307           | 313         | 321          | rBV      | 22809          | 37036         | 1.14%           | 0.310%        |
| 2         | 5.597       | 386           | 393         | 406          | rBV      | 210462         | 389170        | 12.00%          | 3.260%        |
| 3         | 7.183       | 656           | 663         | 677          | rBV2     | 116138         | 256010        | 7.90%           | 2.144%        |
| 4         | 7.547       | 717           | 725         | 741          | rBV      | 433425         | 802599        | 24.76%          | 6.722%        |
| 5         | 8.011       | 794           | 804         | 812          | rBV      | 96180          | 173422        | 5.35%           | 1.453%        |
| 6         | 8.334       | 849           | 859         | 871          | rBV      | 412021         | 751353        | 23.18%          | 6.293%        |
| 7         | 9.174       | 995           | 1002        | 1017         | rVB      | 322351         | 649476        | 20.03%          | 5.440%        |
| 8         | 9.456       | 1040          | 1050        | 1059         | rVB      | 20051          | 34585         | 1.07%           | 0.290%        |
| 9         | 10.813      | 1273          | 1281        | 1288         | rBV2     | 104054         | 219020        | 6.76%           | 1.834%        |
| 10        | 11.277      | 1353          | 1360        | 1368         | rVB      | 43630          | 82890         | 2.56%           | 0.694%        |
| 11        | 11.389      | 1373          | 1379        | 1387         | rVB2     | 39698          | 75521         | 2.33%           | 0.633%        |
| 12        | 13.257      | 1689          | 1697        | 1704         | rBV      | 926445         | 1485704       | 45.83%          | 12.444%       |
| 13        | 14.632      | 1924          | 1931        | 1939         | rBV2     | 191791         | 349275        | 10.77%          | 2.925%        |
| 14        | 16.000      | 2159          | 2164        | 2174         | rBV3     | 58847          | 145504        | 4.49%           | 1.219%        |
| 15        | 16.124      | 2178          | 2185        | 2199         | rVB      | 946103         | 1508192       | 46.52%          | 12.632%       |
| 16        | 17.375      | 2389          | 2398        | 2406         | rBV2     | 302111         | 490018        | 15.11%          | 4.104%        |
| 17        | 19.989      | 2836          | 2843        | 2856         | rBV      | 2339849        | 3241949       | 100.00%         | 27.153%       |
| 18        | 21.640      | 3117          | 3124        | 3133         | rBV      | 337992         | 602919        | 18.60%          | 5.050%        |
| 19        | 23.890      | 3501          | 3507        | 3513         | rBV9     | 17936          | 38679         | 1.19%           | 0.324%        |
| 20        | 24.824      | 3652          | 3666        | 3678         | rBV3     | 192968         | 606140        | 18.70%          | 5.077%        |

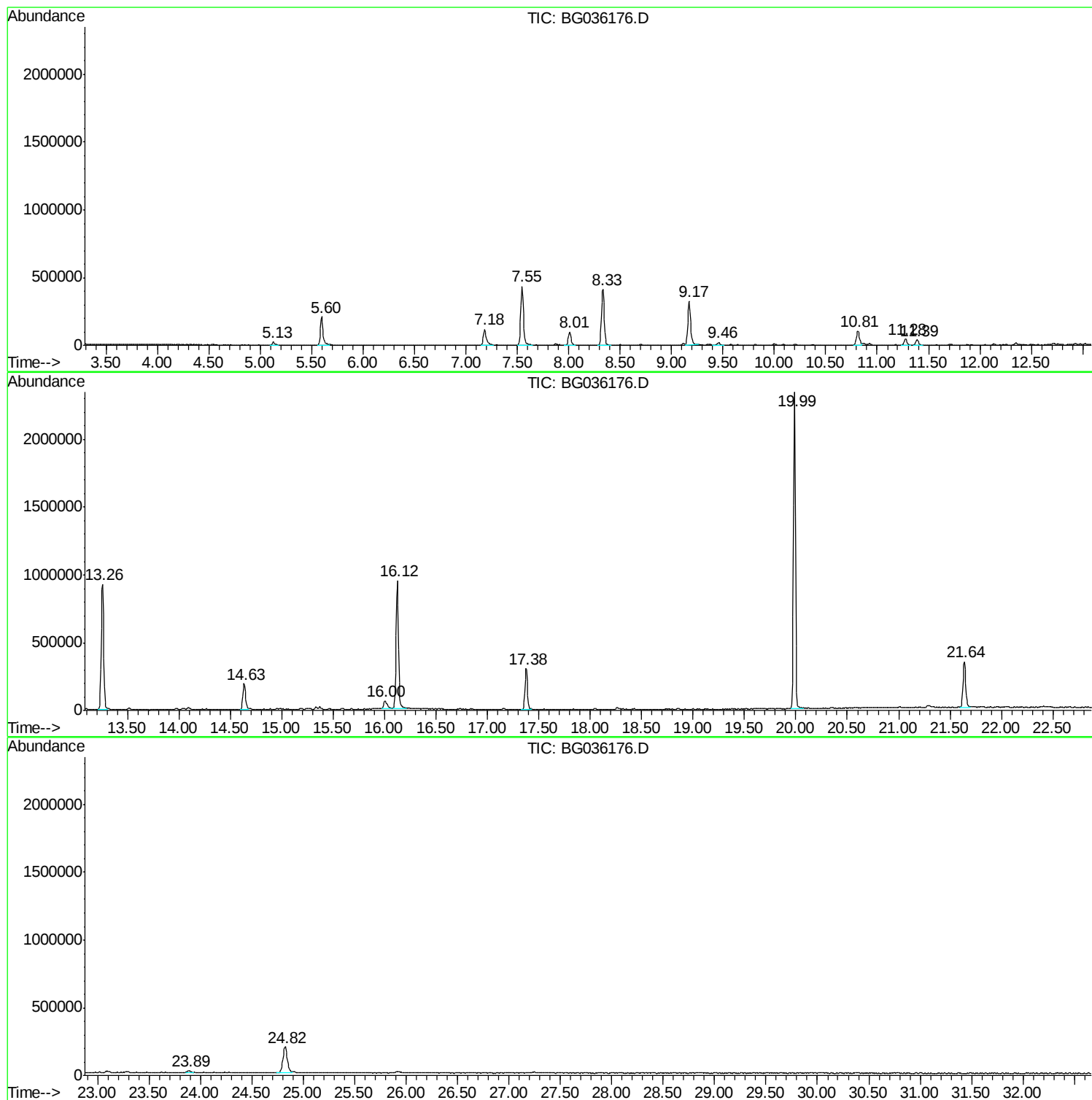
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ALS Vial : 14 Sample Multiplier: 1

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BNA\_G  
ClientSampleId :  
MW-A03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P



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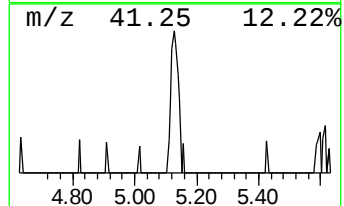
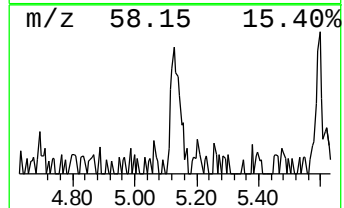
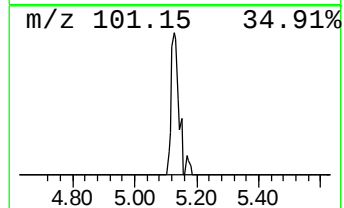
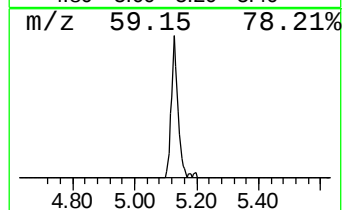
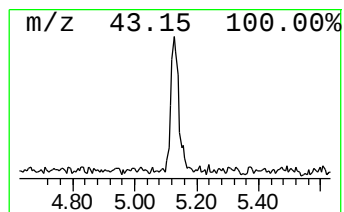
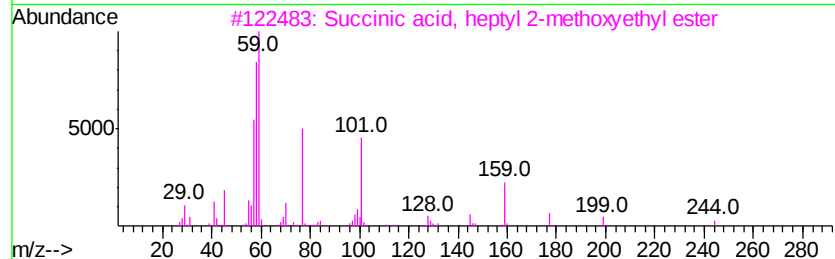
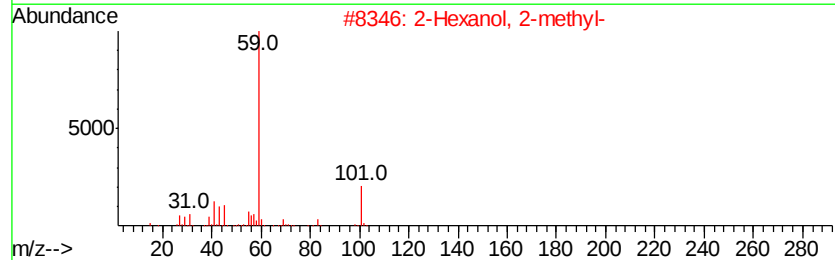
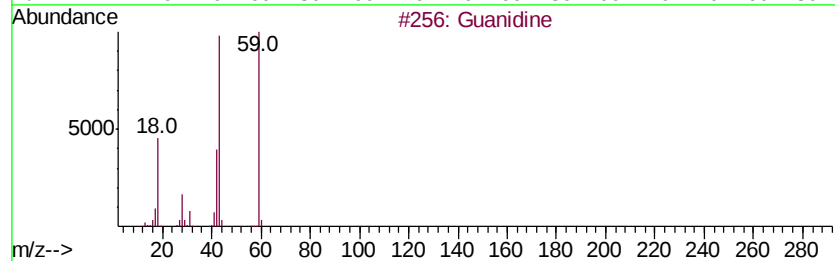
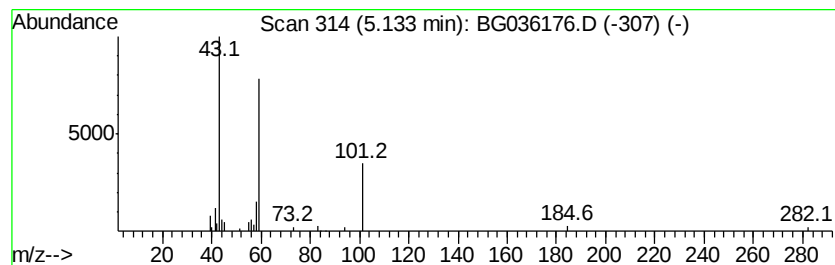
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ClientSampleId :  
MW-A03

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Guanidine Concentration Rank 6

| R.T.      | EstConc                              | Area  | Relative to ISTD       | R.T.         |      |
|-----------|--------------------------------------|-------|------------------------|--------------|------|
| 5.13      | 4.27 ng                              | 37036 | 1,4-Dichlorobenzene-d4 | 8.01         |      |
| Hit# of 5 | Tentative ID                         | MW    | MolForm                | CAS#         | Qual |
| 1         | Guanidine                            | 59    | CH5N3                  | 000113-00-8  | 50   |
| 2         | 2-Hexanol, 2-methyl-                 | 116   | C7H16O                 | 000625-23-0  | 36   |
| 3         | Succinic acid, heptyl 2-methoxyve... | 274   | C14H26O5               | 1000325-80-7 | 36   |
| 4         | 3-Heptanol                           | 116   | C7H16O                 | 000589-82-2  | 33   |
| 5         | Methyl 2,5,8,11-tetraoxatridecan...  | 236   | C10H20O6               | 1000366-78-2 | 33   |



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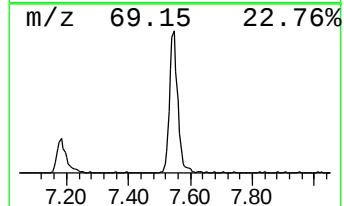
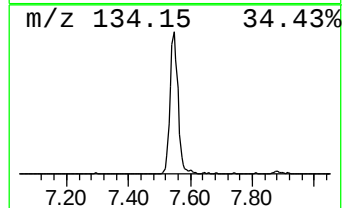
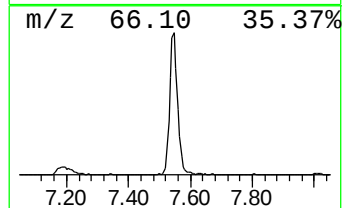
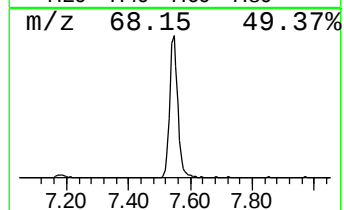
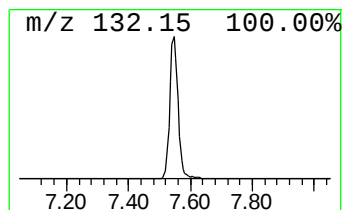
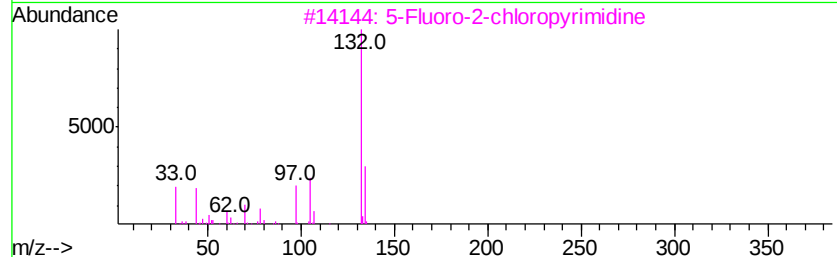
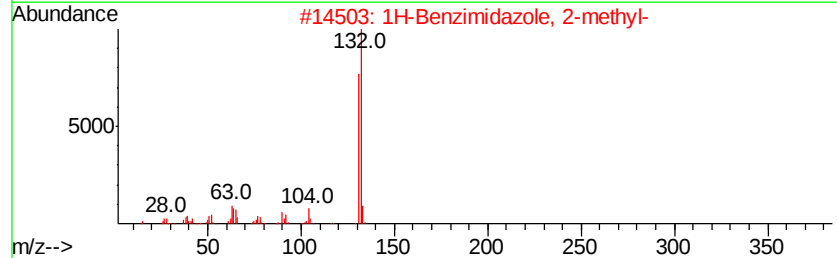
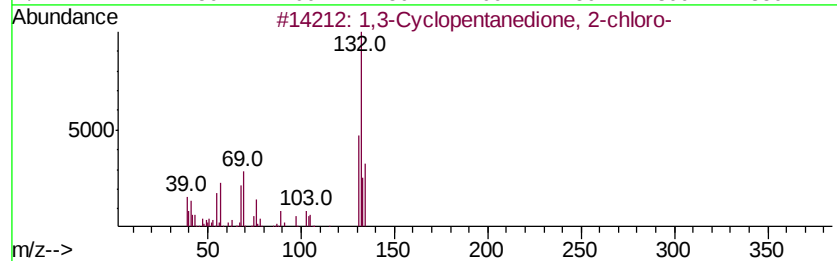
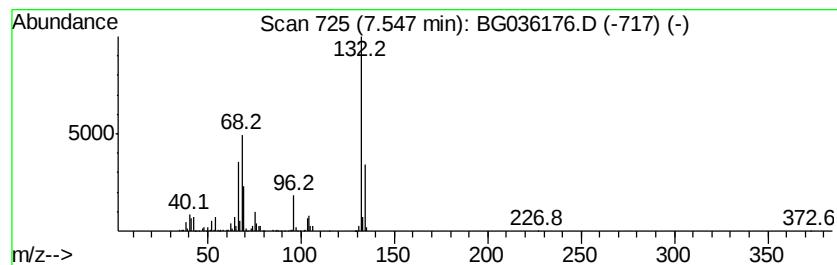
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Peak Number 2 unknown7.55 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.55 | 92.56 ng | 802599 | 1,4-Dichlorobenzene-d4 | 8.01 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 25   |
| 2    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |
| 3    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 4    |    | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 5    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3    | 000372-38-3  | 10   |



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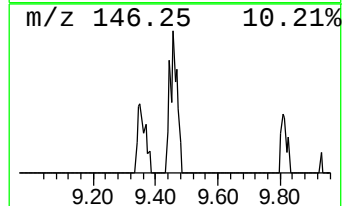
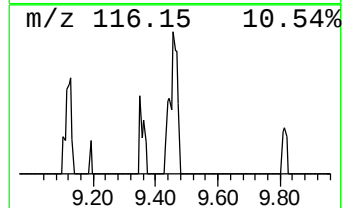
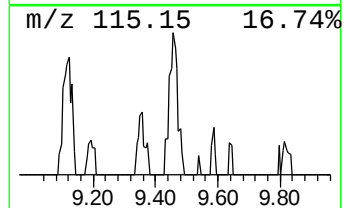
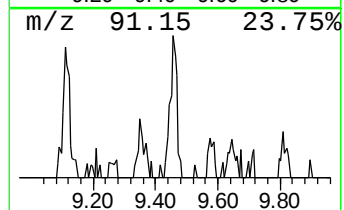
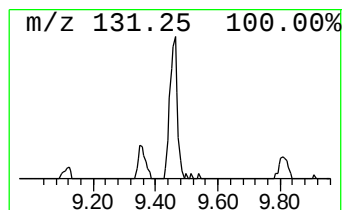
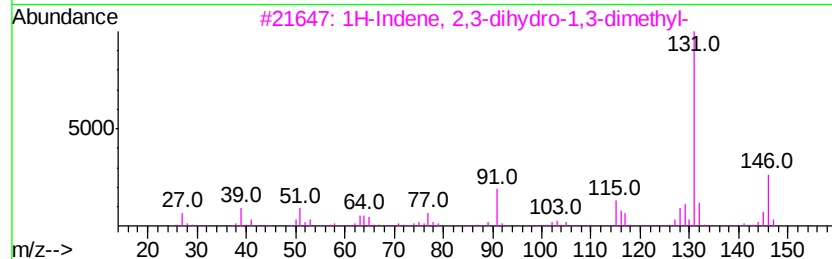
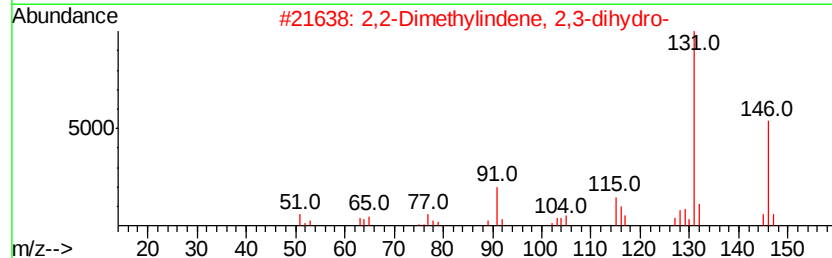
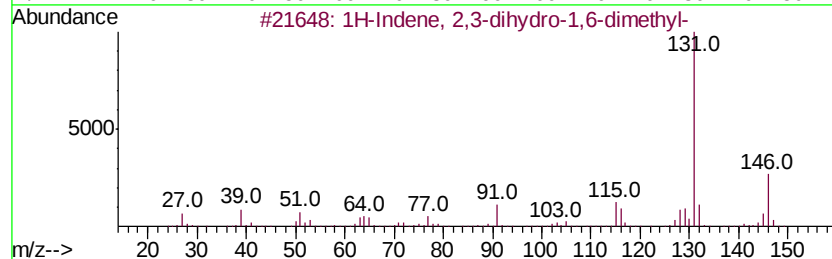
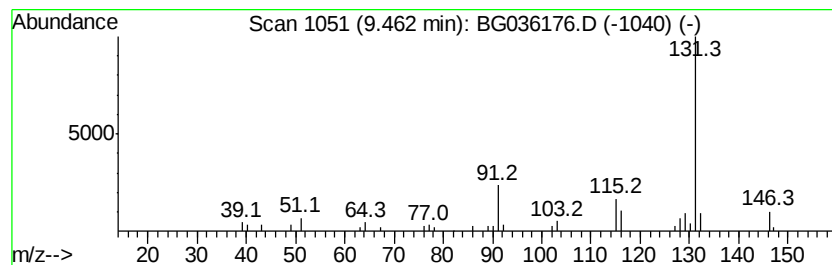
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

| R.T. | EstConc | Area  | Relative to ISTD | R.T.  |
|------|---------|-------|------------------|-------|
| 9.46 | 3.16 ng | 34585 | Napthalene-d8    | 10.81 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14   | 017059-48-2 | 87   |
| 2    |    | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14   | 020836-11-7 | 83   |
| 3    |    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14   | 004175-53-5 | 78   |
| 4    |    | Propanedinitrile, (2,3-dihydro-1... | 196 | C13H12N2 | 069564-96-1 | 72   |
| 5    |    | Benzene, 1-methyl-2-(1-methyl-2-... | 146 | C11H14   | 097664-19-2 | 72   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
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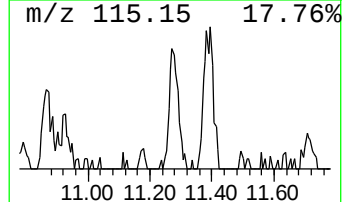
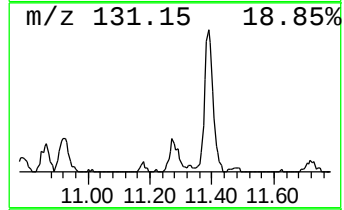
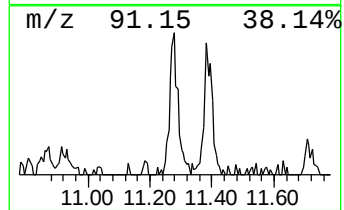
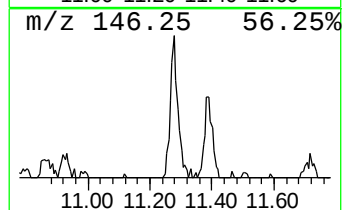
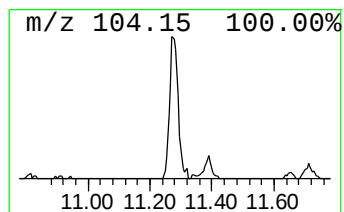
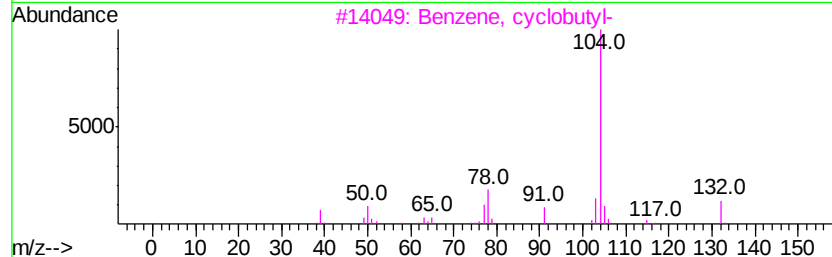
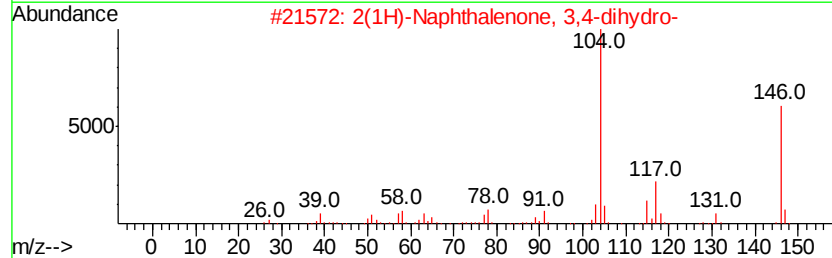
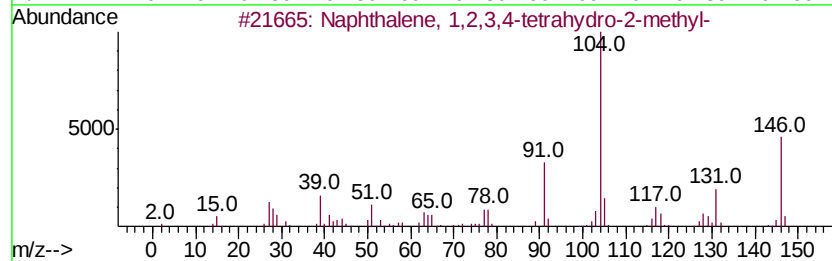
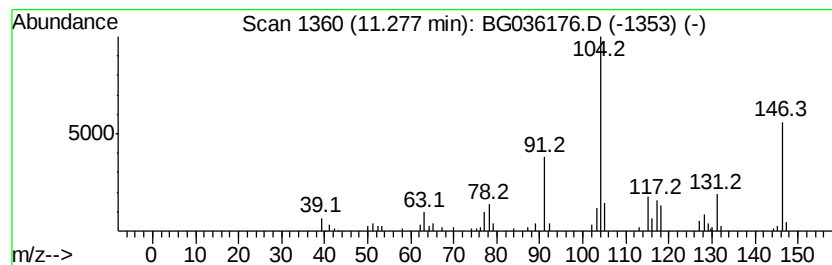
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.28 | 7.57 ng | 82890 | Naphthalene-d8   | 10.81 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------------|-----|----------|-------------|------|
| 1       |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14   | 003877-19-8 | 86   |
| 2       |   | 2(1H)-Naphthalenone, 3,4-dihydro-   | 146 | C10H10O  | 000530-93-8 | 72   |
| 3       |   | Benzene, cyclobutyl-                | 132 | C10H12   | 004392-30-7 | 38   |
| 4       |   | Benzene, 1,1'-(1,2-cyclobutanedi... | 208 | C16H16   | 007694-30-6 | 38   |
| 5       |   | Acetic acid, 2-phenylethyl ester    | 164 | C10H12O2 | 000103-45-7 | 38   |



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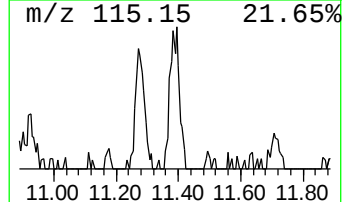
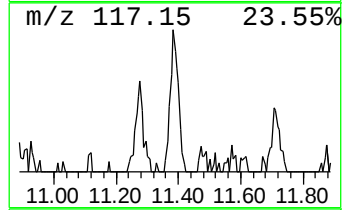
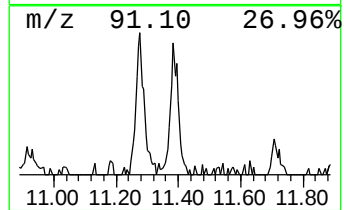
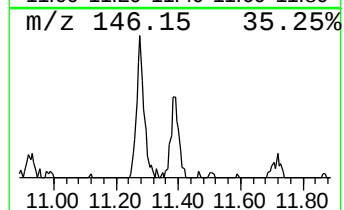
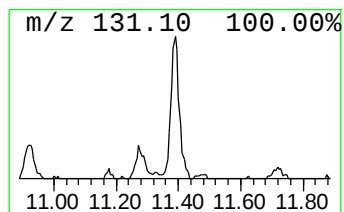
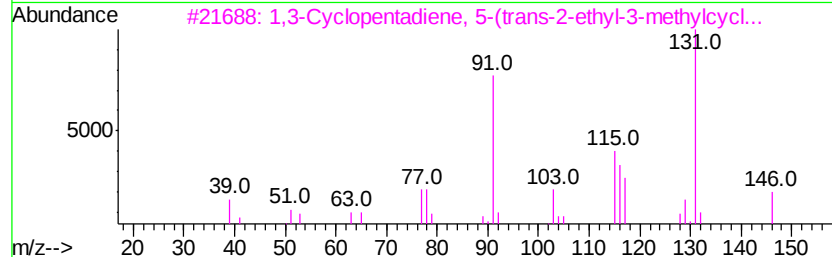
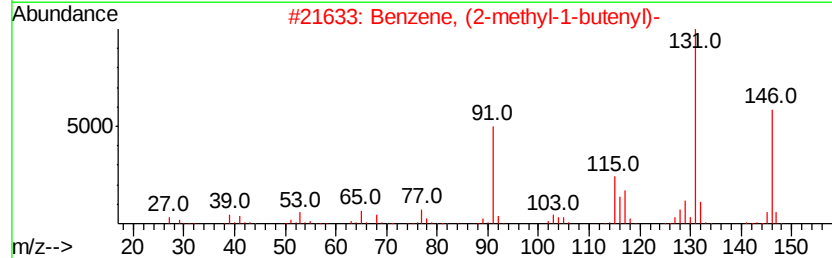
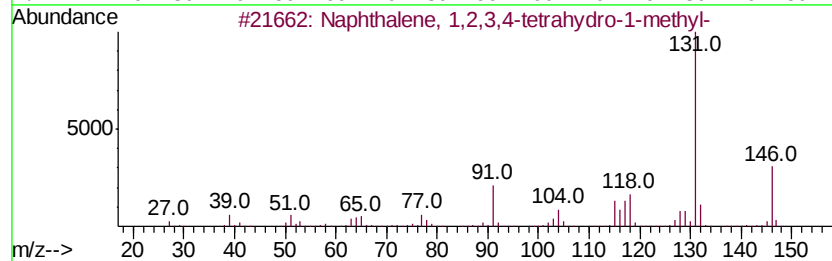
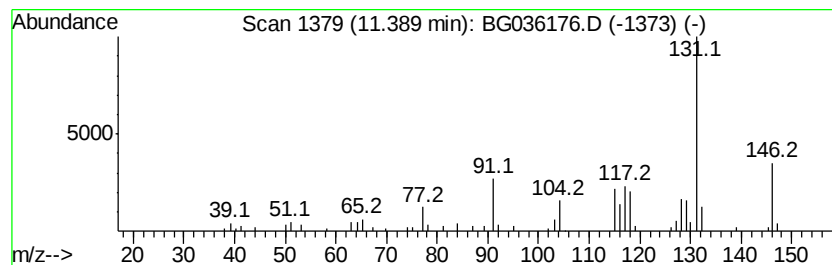
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.39 | 6.90 ng | 75521 | Naphthalene-d8   | 10.81 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------------|-----|---------|-------------|------|
| 1       |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001559-81-5 | 93   |
| 2       |   | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 86   |
| 3       |   | 1,3-Cyclopentadiene, 5-(trans-2-... | 146 | C11H14  | 079209-36-2 | 81   |
| 4       |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 81   |
| 5       |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080718\  
Data File : BG036176.D  
Acq On : 8 Aug 2018 00:18  
Operator : JU/SJ  
Sample : J4303-07  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

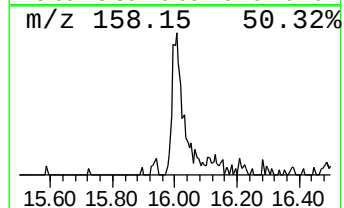
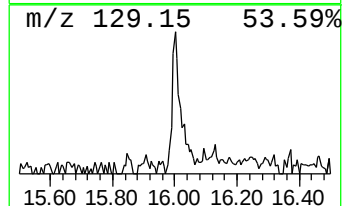
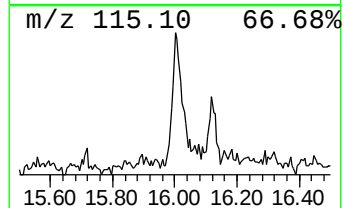
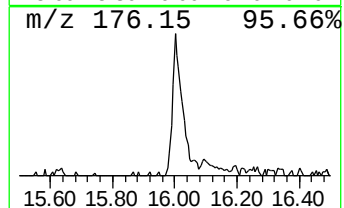
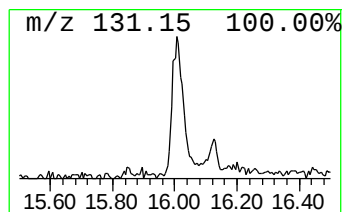
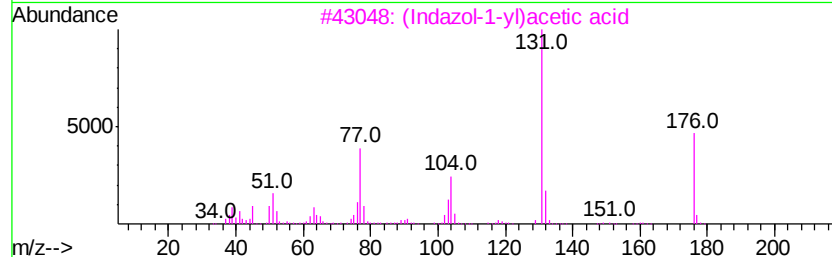
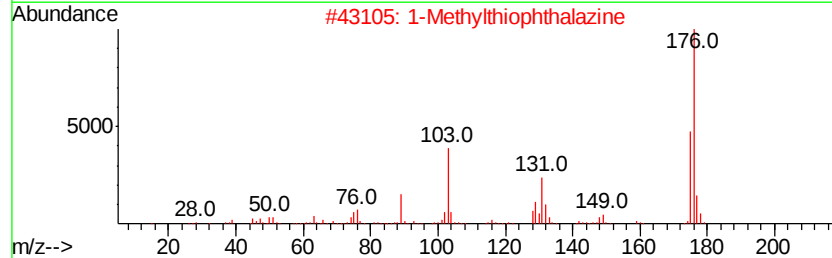
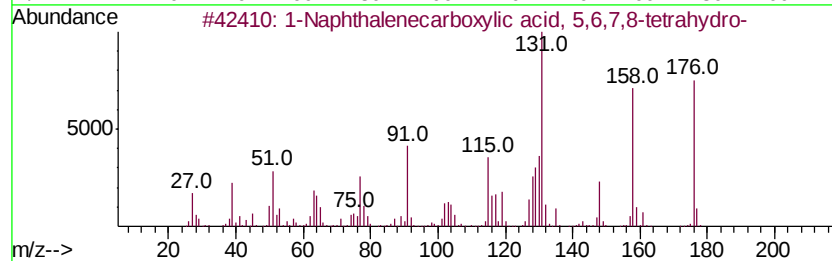
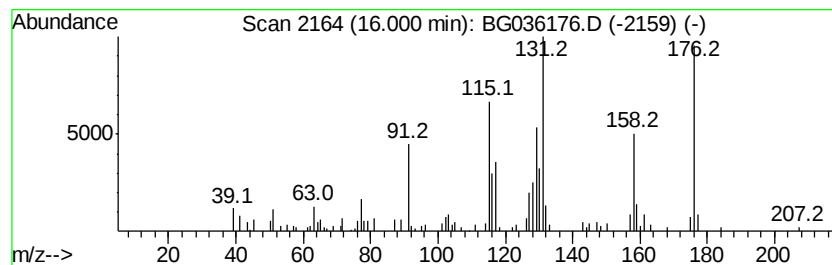
Instrument :  
BNA\_G  
ClientSampleId :  
MW-A03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 1-Naphthalenecarboxylic aci... Concentration Rank 3

| R.T.  | EstConc | Area   | Relative to ISTD                    |     | R.T.     |              |      |
|-------|---------|--------|-------------------------------------|-----|----------|--------------|------|
| 16.00 | 8.33 ng | 145504 | Acenaphthene-d10                    |     | 14.63    |              |      |
| Hit#  | of      | 5      | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
| 1     |         |        | 1-Naphthalenecarboxylic acid, 5,... | 176 | C11H12O2 | 004242-18-6  | 74   |
| 2     |         |        | 1-Methylthiophthalazine             | 176 | C9H8N2S  | 021948-74-3  | 47   |
| 3     |         |        | (Indazol-1-yl)acetic acid           | 176 | C9H8N2O2 | 1000315-94-2 | 42   |
| 4     |         |        | Cinnoline, 3,4-dimethyl-            | 158 | C10H10N2 | 003929-83-7  | 38   |
| 5     |         |        | 1,2-Disilacyclopentane, 1,1,2,2-... | 158 | C7H18Si2 | 015003-82-4  | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG080718\  
Data File : BG036176.D  
Acq On : 8 Aug 2018 00:18  
Operator : JU/SJ  
Sample : J4303-07  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Guanidine            | 5.13  | 4.3     | ng    | 37036    | 1                     | 8.01  | 173422 | 20.0 |
| unknown7.55          | 7.55  | 92.6    | ng    | 802599   | 1                     | 8.01  | 173422 | 20.0 |
| 1H-Indene, 2,3-di... | 9.46  | 3.2     | ng    | 34585    | 2                     | 10.81 | 219020 | 20.0 |
| Naphthalene, 1,2,... | 11.28 | 7.6     | ng    | 82890    | 2                     | 10.81 | 219020 | 20.0 |
| Naphthalene, 1,2,... | 11.39 | 6.9     | ng    | 75521    | 2                     | 10.81 | 219020 | 20.0 |
| 1-Naphthalenecarb... | 16.00 | 8.3     | ng    | 145504   | 3                     | 14.63 | 349275 | 20.0 |