

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
 Data File : BG027364.D  
 Acq On : 10 Jun 2017 2:02  
 Operator : SJ/MA  
 Sample : I3575-06  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 HR-06-060817-B

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:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.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         | 4.551       | 142           | 147         | 156          | rVB      | 84993          | 131858        | 2.81%           | 0.534%        |
| 2         | 5.127       | 239           | 245         | 254          | rBV      | 235490         | 384602        | 8.18%           | 1.559%        |
| 3         | 5.221       | 254           | 261         | 285          | rBV      | 316335         | 681636        | 14.50%          | 2.762%        |
| 4         | 5.703       | 338           | 343         | 355          | rVB      | 41590          | 69865         | 1.49%           | 0.283%        |
| 5         | 6.143       | 410           | 418         | 428          | rBV      | 1105979        | 1893057       | 40.28%          | 7.671%        |
| 6         | 7.642       | 666           | 673         | 677          | rBV      | 500885         | 856101        | 18.22%          | 3.469%        |
| 7         | 7.695       | 677           | 682         | 700          | rVB2     | 840306         | 1624764       | 34.57%          | 6.584%        |
| 8         | 8.088       | 741           | 749         | 763          | rBV      | 1121088        | 2062515       | 43.89%          | 8.358%        |
| 9         | 8.558       | 819           | 829         | 838          | rVB      | 171821         | 321364        | 6.84%           | 1.302%        |
| 10        | 8.881       | 874           | 884         | 894          | rBV      | 724098         | 1381905       | 29.41%          | 5.600%        |
| 11        | 9.111       | 914           | 923         | 932          | rBV3     | 56509          | 107428        | 2.29%           | 0.435%        |
| 12        | 9.716       | 1018          | 1026        | 1036         | rBV      | 671087         | 1291881       | 27.49%          | 5.235%        |
| 13        | 10.086      | 1081          | 1089        | 1099         | rBV      | 113865         | 186847        | 3.98%           | 0.757%        |
| 14        | 11.373      | 1295          | 1308        | 1316         | rBV      | 270576         | 524474        | 11.16%          | 2.125%        |
| 15        | 12.565      | 1505          | 1511        | 1519         | rBV3     | 61557          | 132436        | 2.82%           | 0.537%        |
| 16        | 13.764      | 1707          | 1715        | 1723         | rBV      | 1964668        | 3177002       | 67.60%          | 12.874%       |
| 17        | 14.069      | 1761          | 1767        | 1774         | rBV2     | 41615          | 104586        | 2.23%           | 0.424%        |
| 18        | 15.139      | 1942          | 1949        | 1958         | rVB      | 279781         | 497257        | 10.58%          | 2.015%        |
| 19        | 16.619      | 2193          | 2201        | 2222         | rBV      | 1692919        | 2722027       | 57.92%          | 11.031%       |
| 20        | 17.883      | 2407          | 2416        | 2430         | rBV2     | 538683         | 916466        | 19.50%          | 3.714%        |
| 21        | 20.462      | 2848          | 2855        | 2862         | rBV      | 3475327        | 4699507       | 100.00%         | 19.044%       |
| 22        | 22.266      | 3154          | 3162        | 3175         | rBV2     | 336308         | 748187        | 15.92%          | 3.032%        |
| 23        | 25.956      | 3783          | 3790        | 3804         | rVB6     | 45144          | 161120        | 3.43%           | 0.653%        |

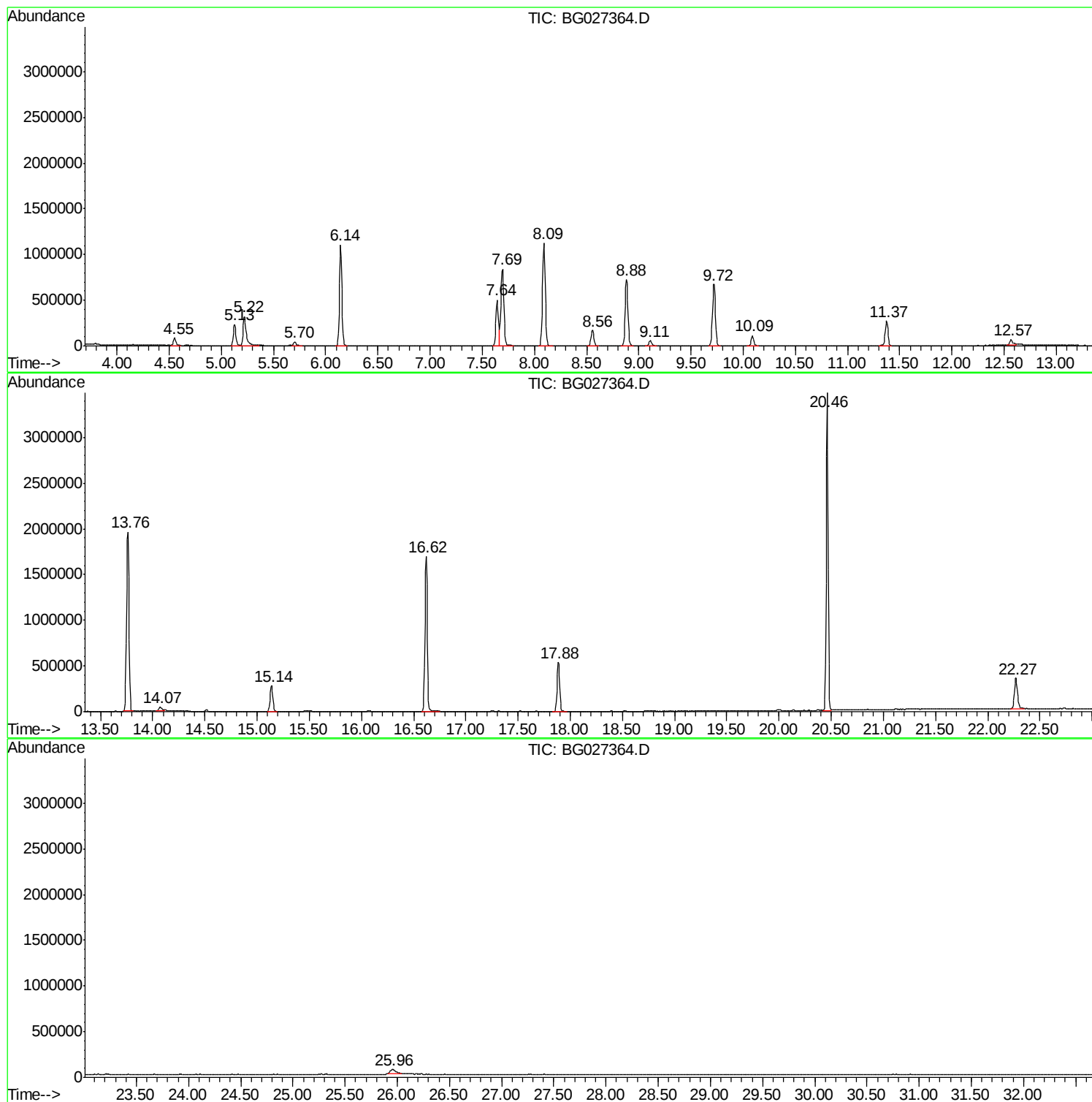
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ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.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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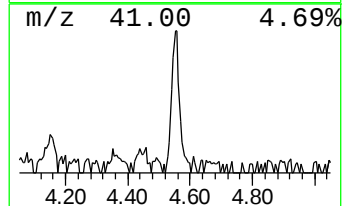
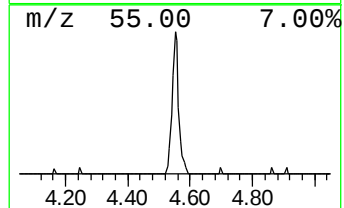
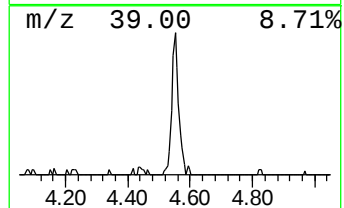
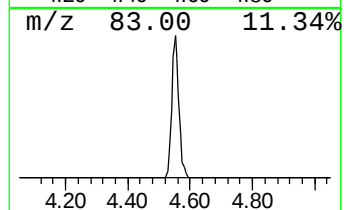
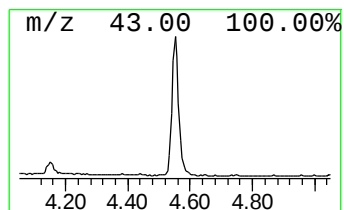
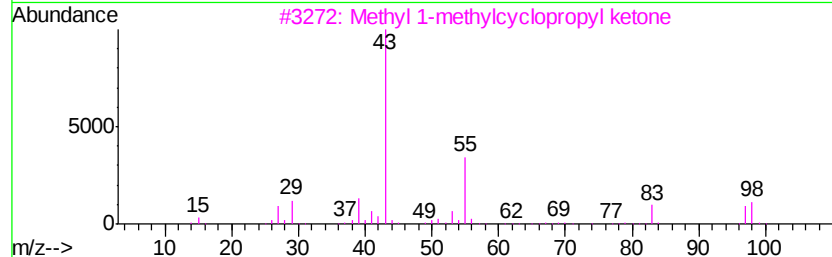
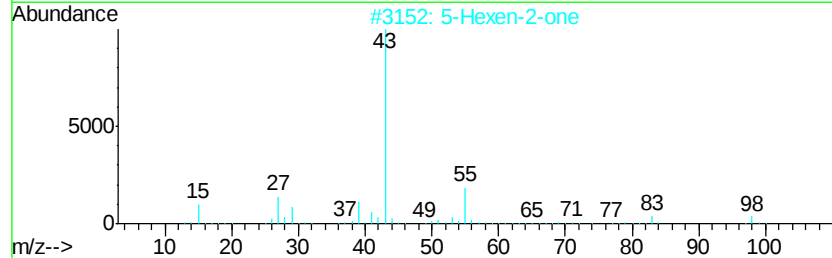
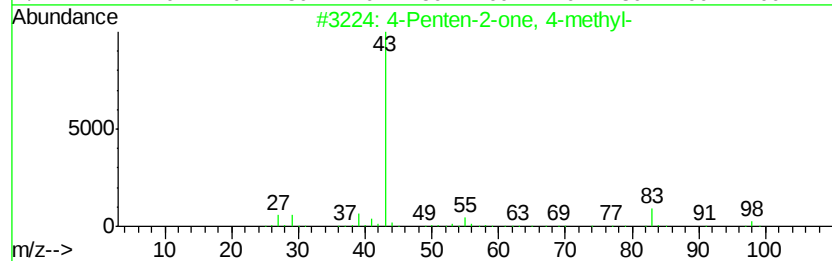
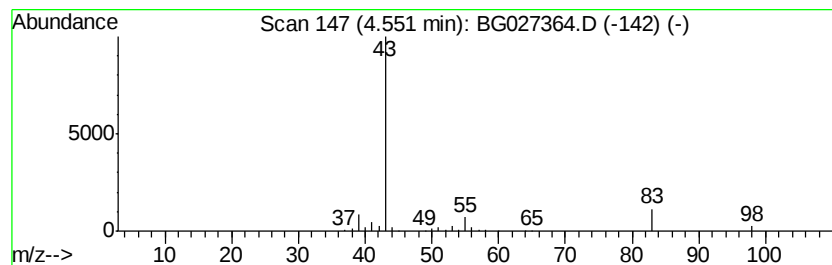
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.55 | 8.21 ng | 131858 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4-Penten-2-one, 4-methyl-         | 98  | C6H10O  | 003744-02-3 | 72   |
| 2    |    | 5-Hexen-2-one                     | 98  | C6H10O  | 000109-49-9 | 35   |
| 3    |    | Methyl 1-methylcyclopropyl ketone | 98  | C6H10O  | 001567-75-5 | 30   |
| 4    |    | 2-Vinylethyl acetate              | 114 | C6H10O2 | 001576-84-7 | 28   |
| 5    |    | 2-Pentanone, 3-methylene-         | 98  | C6H10O  | 004359-77-7 | 17   |



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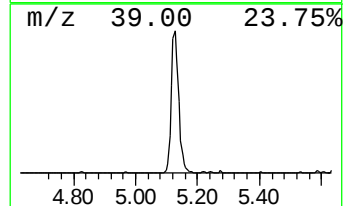
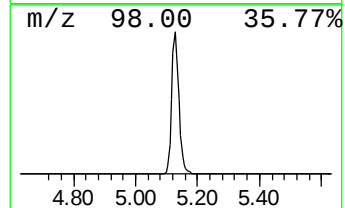
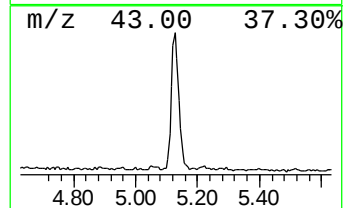
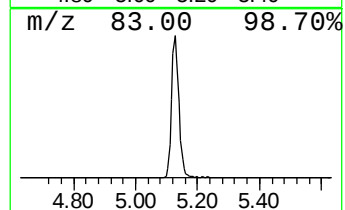
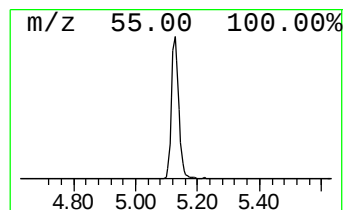
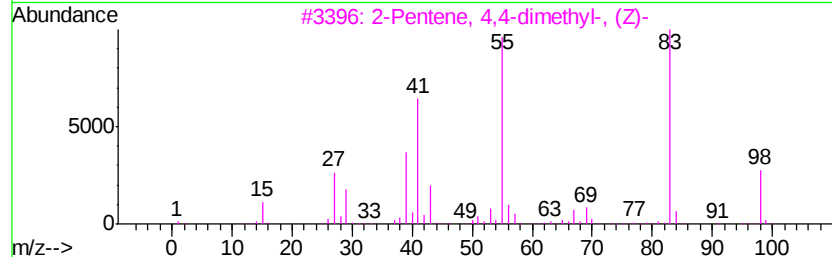
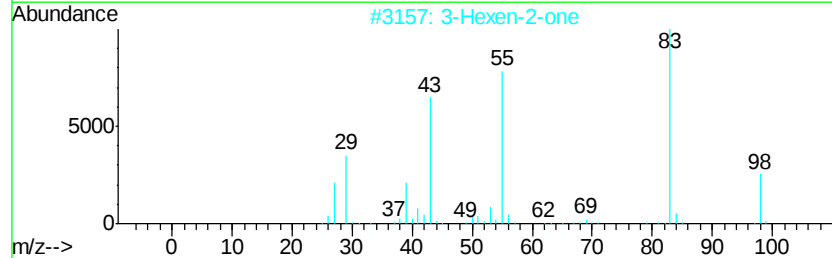
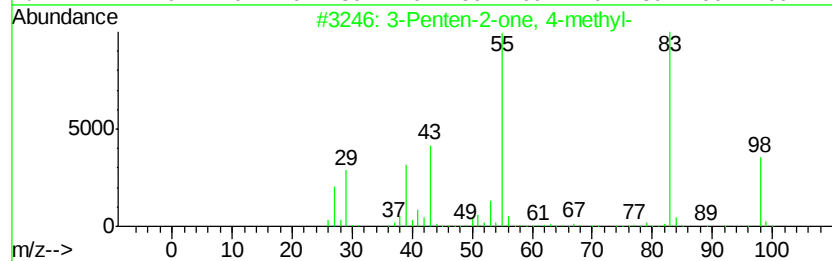
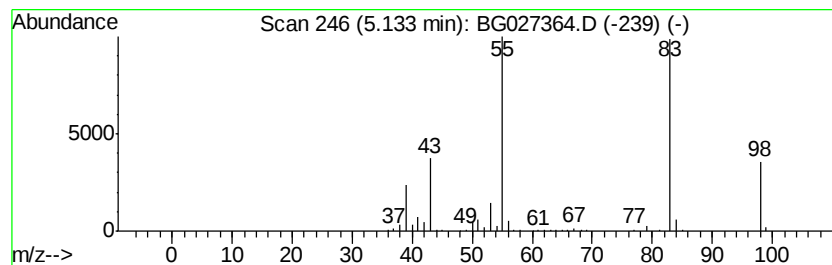
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.13 | 23.94 ng | 384602 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|--------------------------------|----|---------|-------------|------|
| 1    | 5  | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 94   |
| 2    |    | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 91   |
| 3    |    | 2-Pentene, 4,4-dimethyl-, (Z)- | 98 | C7H14   | 000762-63-0 | 86   |
| 4    |    | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 86   |
| 5    |    | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 86   |



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BNA\_G  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

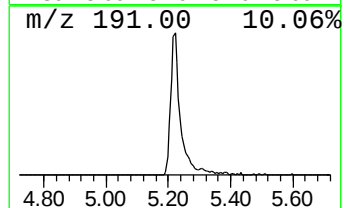
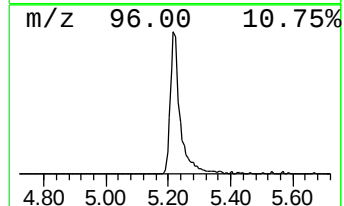
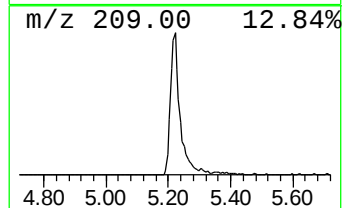
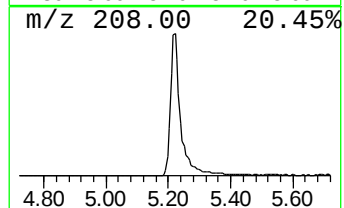
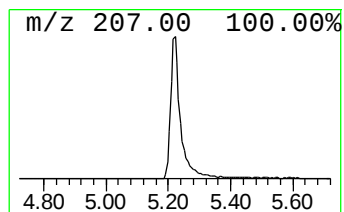
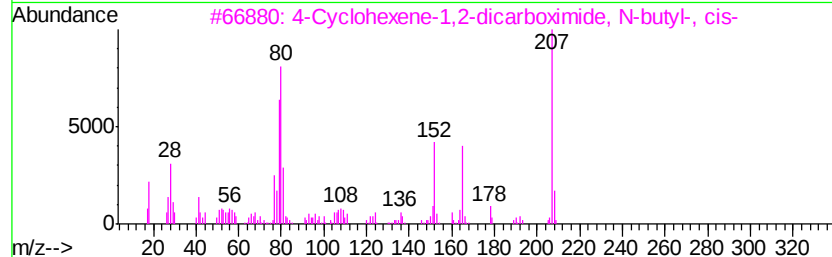
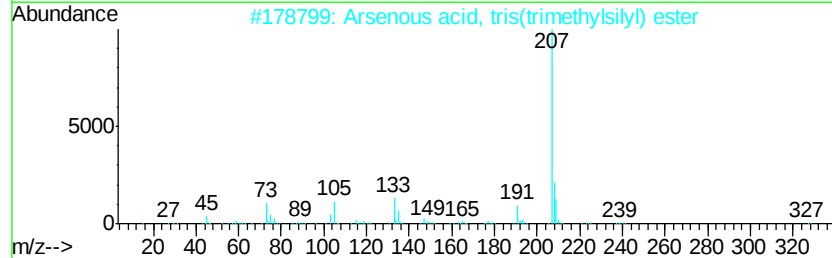
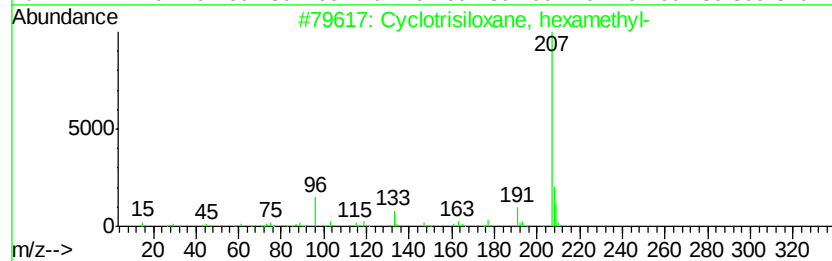
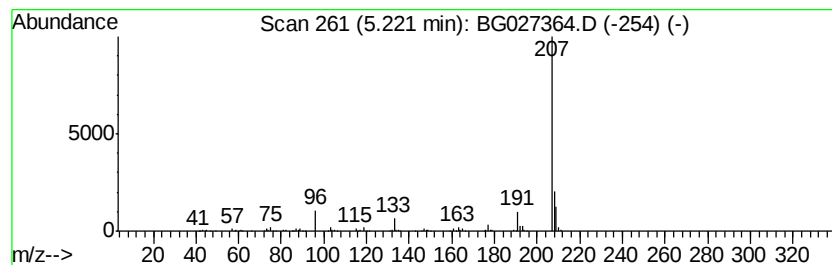
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Cyclotrisiloxane, hexamethyl- Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.22 | 42.42 ng | 681636 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#        | Qual |
|------|----|-------------------------------------|-----|--------------|-------------|------|
| 1    | 5  | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3   | 000541-05-9 | 91   |
| 2    |    | Arsenous acid, tris(trimethylsil... | 342 | C9H27AsO3Si3 | 055429-29-3 | 64   |
| 3    |    | 4-Cyclohexene-1,2-dicarboximide,... | 207 | C12H17N02    | 028916-00-9 | 42   |
| 4    |    | 1H-Indole, 1-methyl-2-phenyl-       | 207 | C15H13N      | 003558-24-5 | 39   |
| 5    |    | Benzo[h]quinoline, 2,4-dimethyl-    | 207 | C15H13N      | 000605-67-4 | 38   |



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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

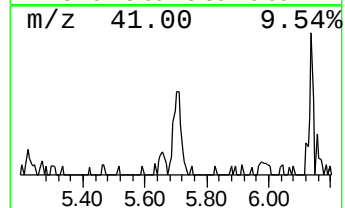
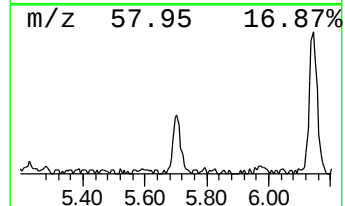
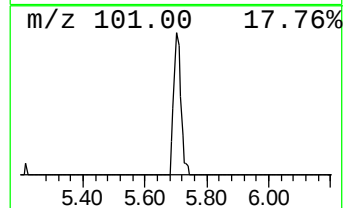
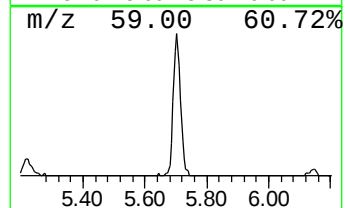
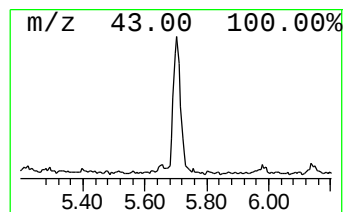
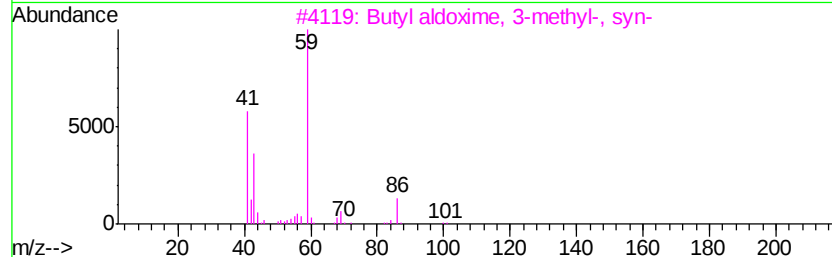
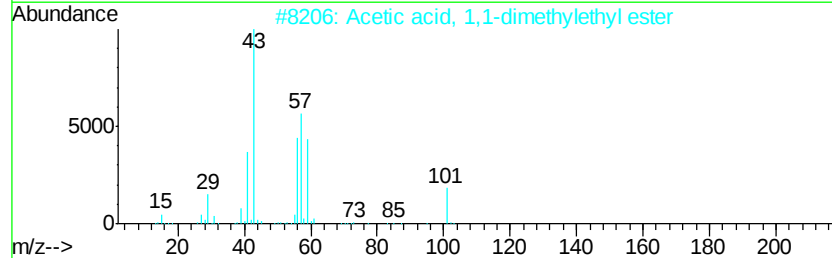
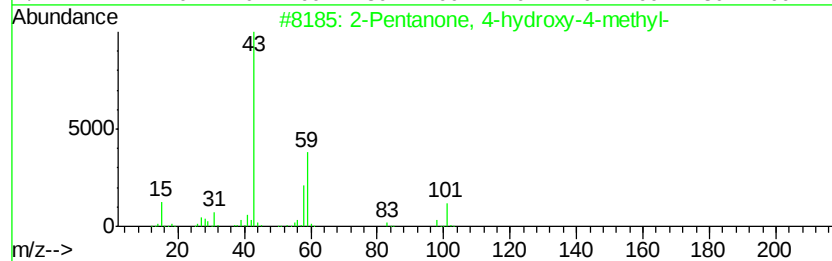
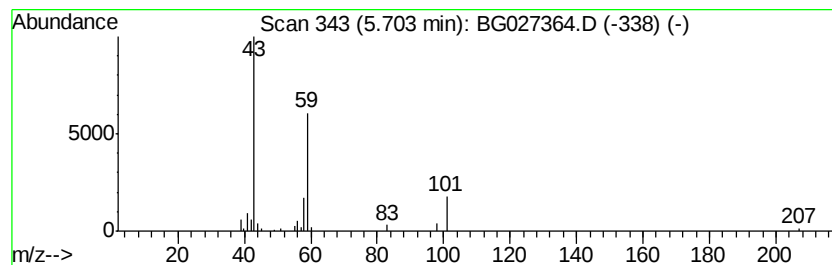
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.70 | 4.35 ng | 69865 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 38   |
| 3    |    | Butyl aldoxime, 3-methyl-, syn-     | 101 | C5H11NO  | 005780-40-5 | 23   |
| 4    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 5    |    | Acetone                             | 58  | C3H6O    | 000067-64-1 | 9    |



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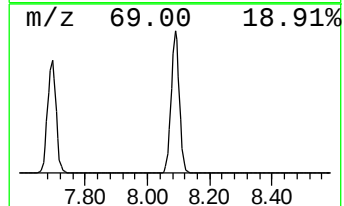
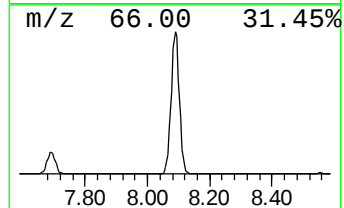
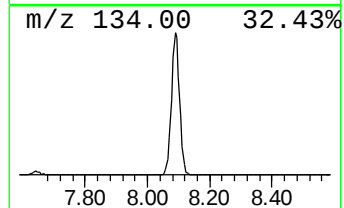
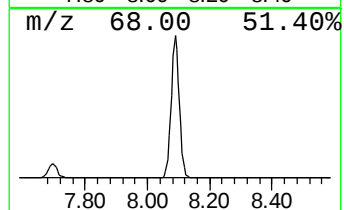
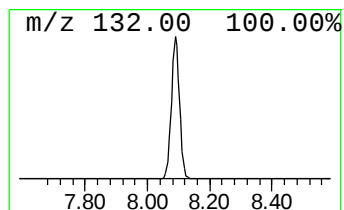
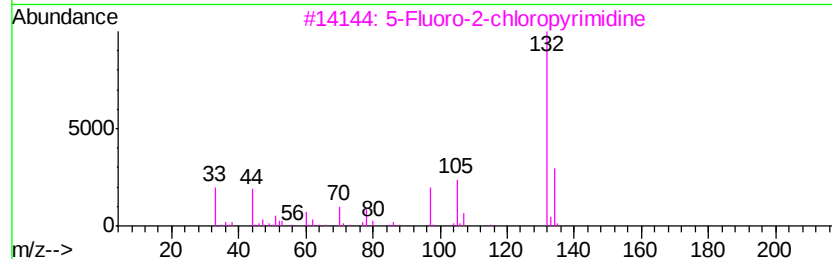
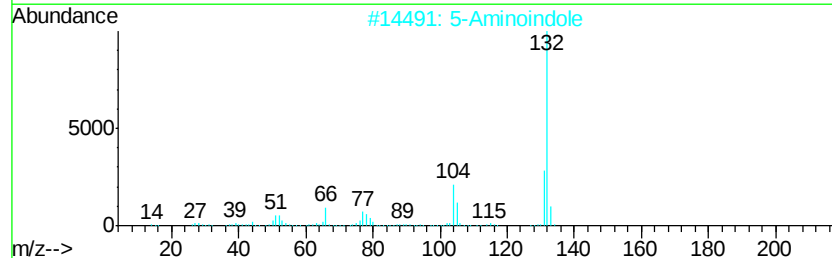
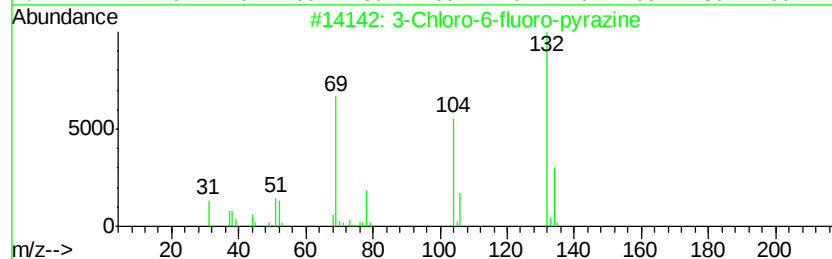
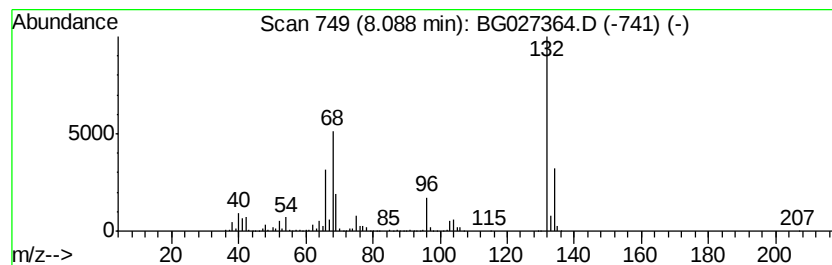
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown8.09 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 8.09 | 128.36 ng | 2062520 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of                          | Tentative ID | MW        | MolForm      | CAS# | Qual |
|------|-----------------------------|--------------|-----------|--------------|------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine  | 132          | C4H2ClFN2 | 1000146-10-7 | 16   |      |
| 2    | 5-Aminoindole               | 132          | C8H8N2    | 005192-03-0  | 12   |      |
| 3    | 5-Fluoro-2-chloropyrimidine | 132          | C4H2ClFN2 | 062802-42-0  | 12   |      |
| 4    | 2-Cyclohexen-1-one          | 96           | C6H8O     | 000930-68-7  | 10   |      |
| 5    | 2-Propyn-1-amine, N-methyl- | 69           | C4H7N     | 035161-71-8  | 9    |      |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027364.D  
Acq On : 10 Jun 2017 2:02  
Operator : SJ/MA  
Sample : I3575-06  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

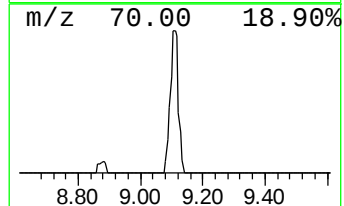
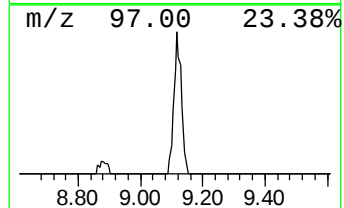
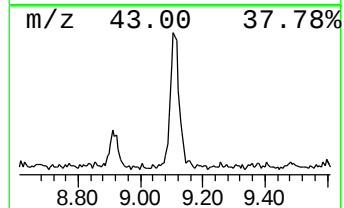
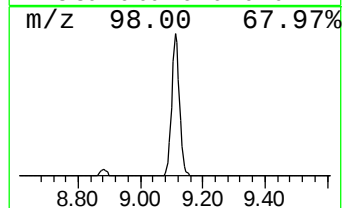
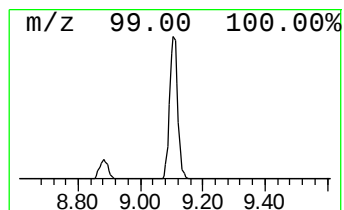
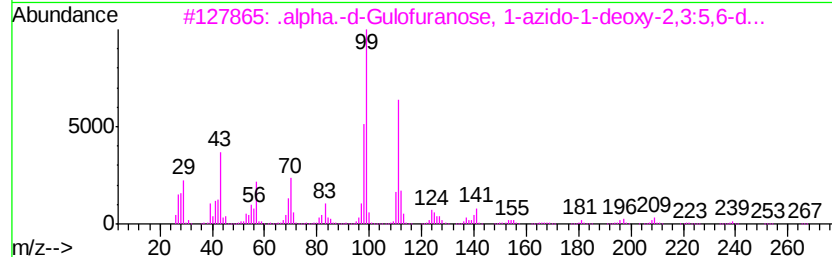
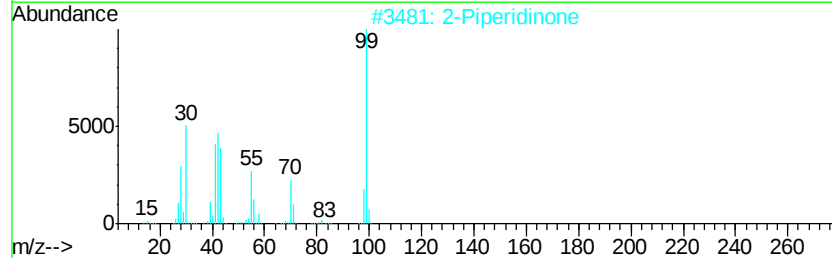
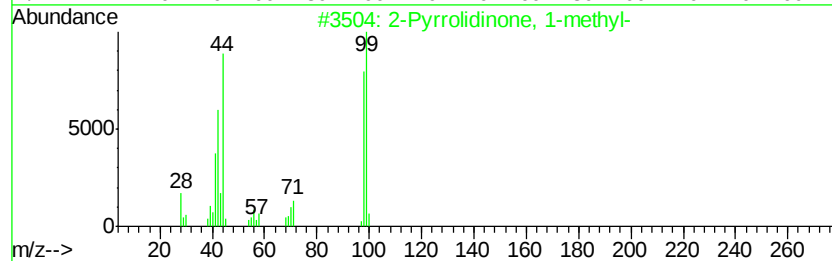
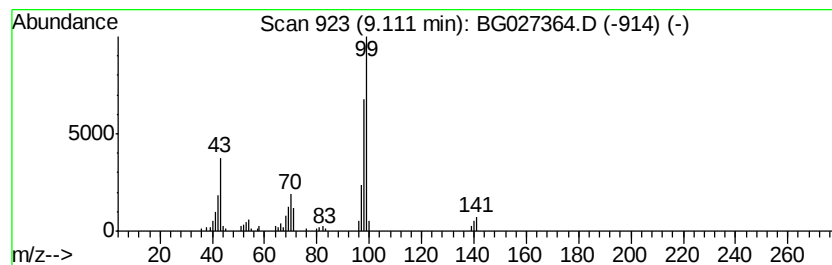
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 2-Pyrrolidinone, 1-methyl- Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 9.11 | 6.69 ng | 107428 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|---------|---|-------------------------------------|-----|--------------|--------------|------|
| 1       |   | 2-Pyrrolidinone, 1-methyl-          | 99  | C5H9NO       | 000872-50-4  | 64   |
| 2       |   | 2-Piperidinone                      | 99  | C5H9NO       | 000675-20-7  | 49   |
| 3       |   | .alpha.-d-Gulofuranose, 1-azido-... | 281 | C10H17B2N3O5 | 1000149-82-2 | 39   |
| 4       |   | n-Heptyl methylphosphonofluoridate  | 196 | C8H18FO2P    | 162085-82-7  | 37   |
| 5       |   | Imidosulfurous difluoride, methyl-  | 99  | CH3F2NS      | 000758-20-3  | 36   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027364.D  
Acq On : 10 Jun 2017 2:02  
Operator : SJ/MA  
Sample : I3575-06  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

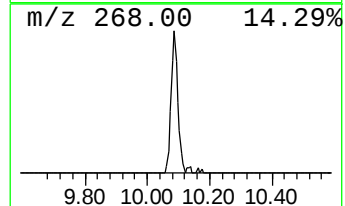
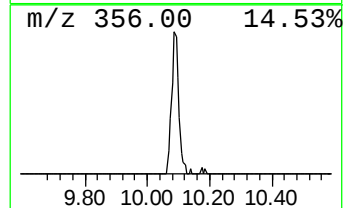
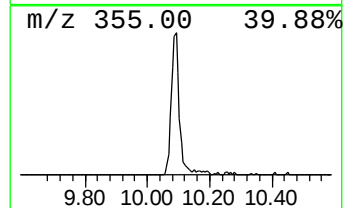
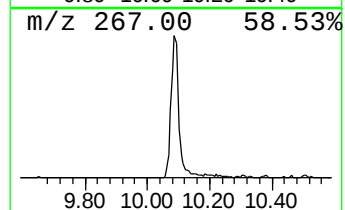
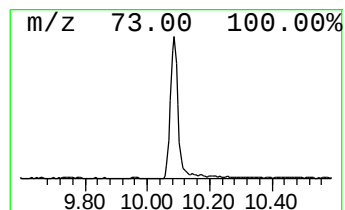
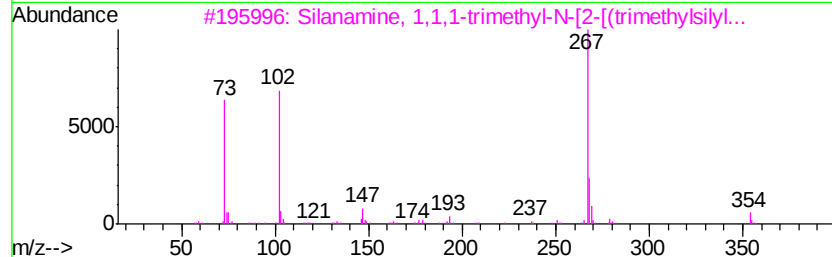
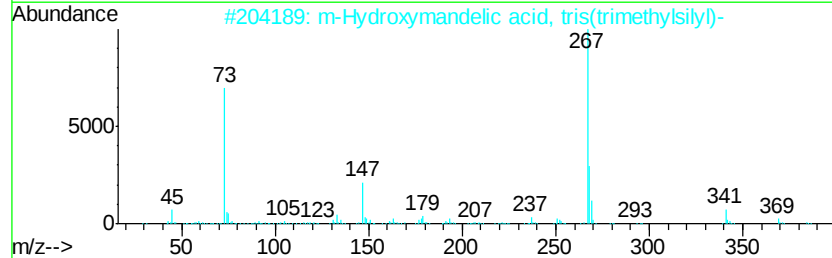
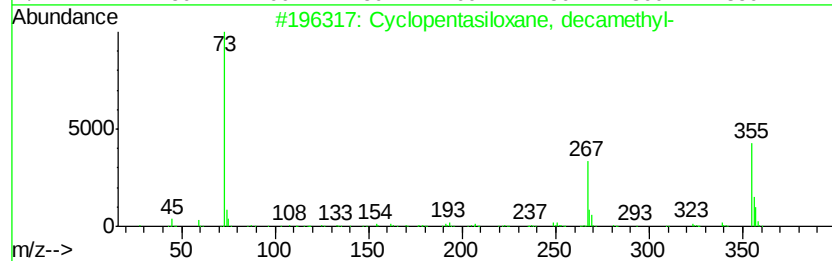
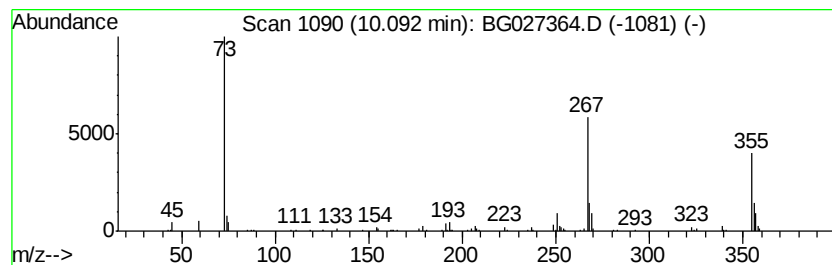
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Cyclopentasiloxane, decamet... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.09 | 7.13 ng | 186847 | Naphthalene-d8   | 11.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#        | Qual |
|------|----|-------------------------------------|-----|--------------|-------------|------|
| 1    | 5  | Cyclopentasiloxane, decamethyl-     | 370 | C10H30O5Si5  | 000541-02-6 | 91   |
| 2    |    | m-Hydroxymandelic acid, tris(tri... | 384 | C17H32O4Si3  | 068595-69-7 | 50   |
| 3    |    | Silanamine, 1,1,1-trimethyl-N-[2... | 369 | C17H35N02Si3 | 068595-60-8 | 50   |
| 4    |    | Benzoic acid, 2-[(trimethylsilyl... | 282 | C13H22O3Si2  | 003789-85-3 | 47   |
| 5    |    | Benzeneethanamine, N-butyl-.beta... | 353 | C18H35N02Si2 | 040629-66-1 | 47   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027364.D  
Acq On : 10 Jun 2017 2:02  
Operator : SJ/MA  
Sample : I3575-06  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

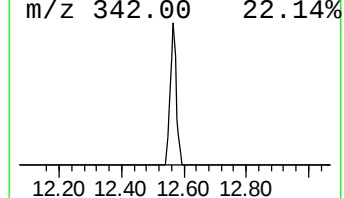
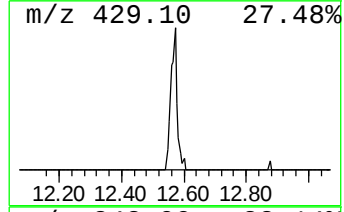
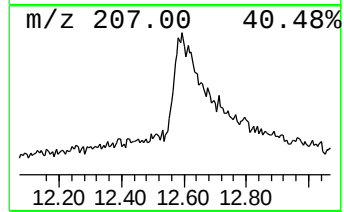
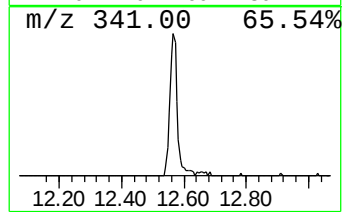
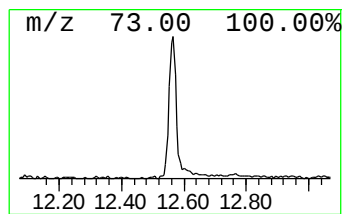
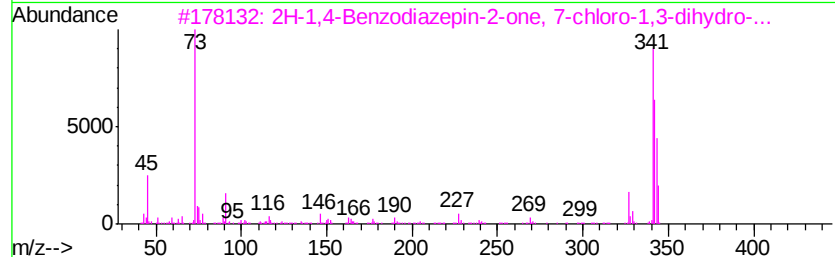
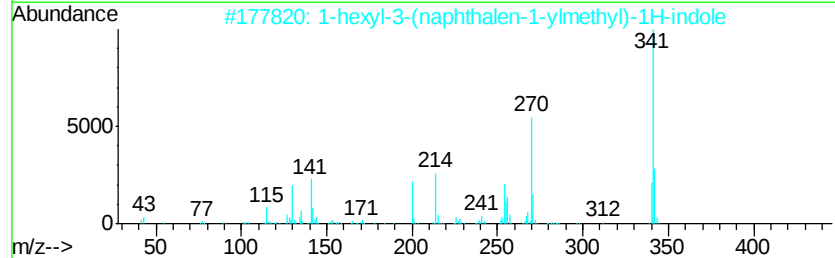
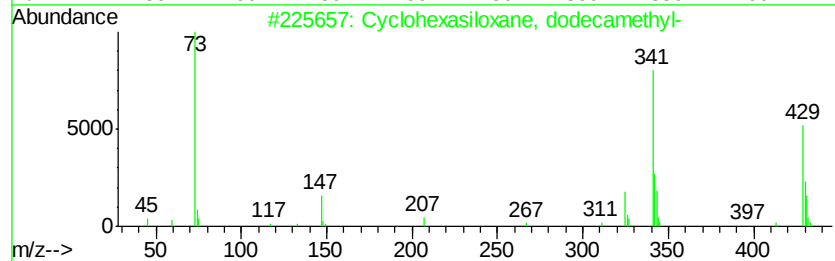
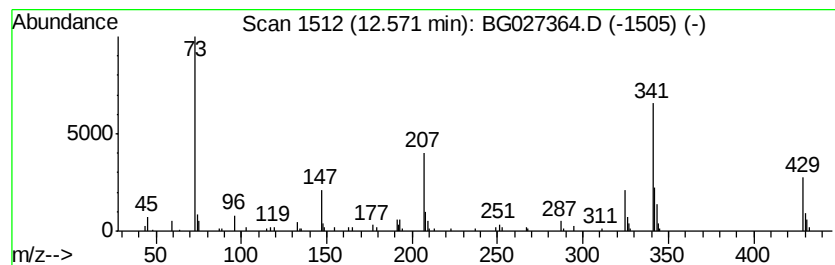
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Cyclohexasiloxane, dodecane... Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.57 | 5.05 ng | 132436 | Naphthalene-d8   | 11.37 |

| Hit# | of | Tentative ID                         | MW  | MolForm       | CAS#         | Qual |
|------|----|--------------------------------------|-----|---------------|--------------|------|
| 1    | 5  | Cyclohexasiloxane, dodecamethyl-     | 444 | C12H36O6Si6   | 000540-97-6  | 91   |
| 2    |    | 1-hexyl-3-(naphthalen-1-ylmethyl)... | 341 | C25H27N       | 1000372-24-4 | 38   |
| 3    |    | 2H-1,4-Benzodiazepin-2-one, 7-ch...  | 342 | C18H19ClN2OSi | 055299-24-6  | 37   |
| 4    |    | 1,3,5,7,9-Pentaethylcyclopentasi...  | 370 | C10H30O5Si5   | 017995-44-7  | 37   |
| 5    |    | Fluoren-9-ol, 3,6-dimethoxy-9-(2...  | 342 | C23H18O3      | 1000217-31-2 | 35   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027364.D  
Acq On : 10 Jun 2017 2:02  
Operator : SJ/MA  
Sample : I3575-06  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

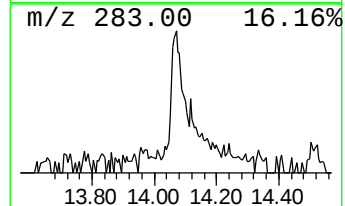
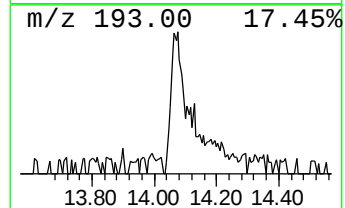
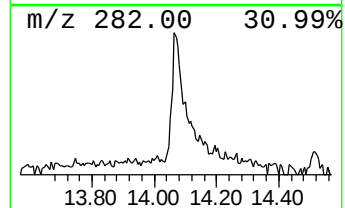
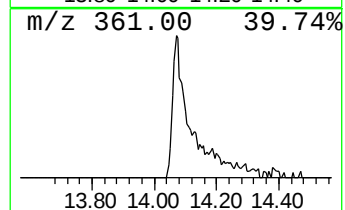
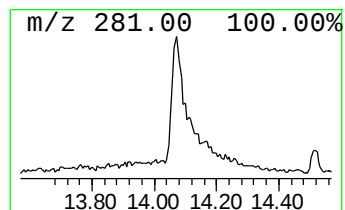
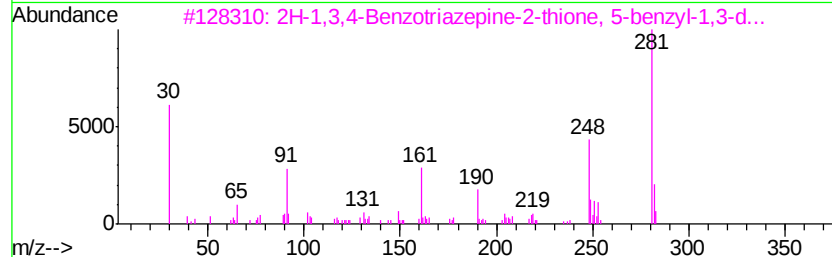
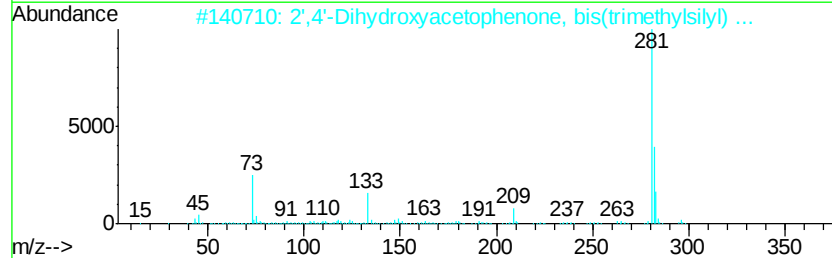
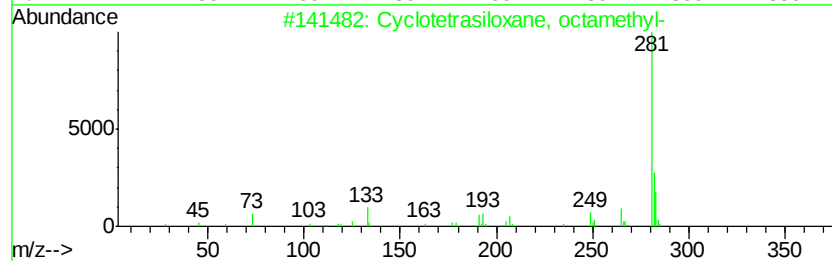
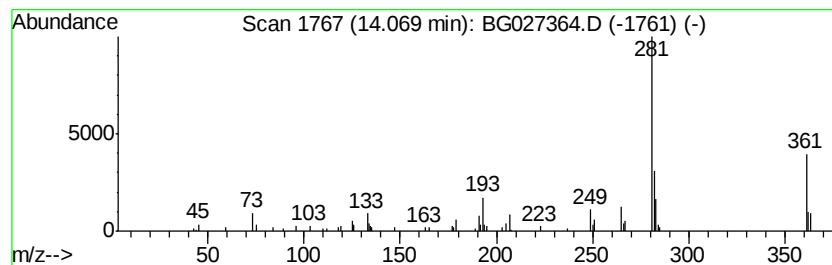
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Cyclotetrasiloxane, octamet... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.07 | 4.21 ng | 104586 | Acenaphthene-d10 | 15.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Cyclotetrasiloxane, octamethyl-     | 296 | C8H24O4Si4   | 000556-67-2  | 52   |
| 2    |    | 2',4'-Dihydroxyacetophenone, bis... | 296 | C14H24O3Si2  | 1000352-82-1 | 43   |
| 3    |    | 2H-1,3,4-Benzotriazepine-2-thion... | 281 | C16H15N3S    | 141071-26-3  | 43   |
| 4    |    | Ethylphosphonic acid, bis(tert-b... | 338 | C14H35O3PSi2 | 1000273-30-1 | 43   |
| 5    |    | 4H-1,2,4-Triazole-3-thiol, 4-all... | 281 | C16H15N3S    | 031803-13-1  | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027364.D  
Acq On : 10 Jun 2017 2:02  
Operator : SJ/MA  
Sample : I3575-06  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HR-06-060817-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.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 |
| 4-Penten-2-one, 4... | 4.55  | 8.2     | ng    | 131858   | 1                     | 8.56  | 321364 | 20.0 |
| 3-Penten-2-one, 4... | 5.13  | 23.9    | ng    | 384602   | 1                     | 8.56  | 321364 | 20.0 |
| Cyclotrisiloxane,... | 5.22  | 42.4    | ng    | 681636   | 1                     | 8.56  | 321364 | 20.0 |
| 2-Pentanone, 4-hy... | 5.70  | 4.3     | ng    | 69865    | 1                     | 8.56  | 321364 | 20.0 |
| unknown8.09          | 8.09  | 128.4   | ng    | 2062520  | 1                     | 8.56  | 321364 | 20.0 |
| 2-Pyrrolidinone, ... | 9.11  | 6.7     | ng    | 107428   | 1                     | 8.56  | 321364 | 20.0 |
| Cyclopentasiloxan... | 10.09 | 7.1     | ng    | 186847   | 2                     | 11.37 | 524474 | 20.0 |
| Cyclohexasiloxane... | 12.57 | 5.0     | ng    | 132436   | 2                     | 11.37 | 524474 | 20.0 |
| Cyclotetrasiloxan... | 14.07 | 4.2     | ng    | 104586   | 3                     | 15.13 | 497257 | 20.0 |