

Data Path : Z:\HPCHEM1\BNA G\DATA\BG052715\  
 Data File : BG017319.D  
 Acq On : 28 May 2015 3:57  
 Operator : TP/IZ  
 Sample : G2387-05  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 TP-16-D8

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

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

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG051315.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         | 2.820       | 18            | 22          | 27           | rVV      | 19833          | 25990         | 1.33%           | 0.238%        |
| 2         | 2.867       | 27            | 30          | 38           | rVB3     | 13864          | 19521         | 1.00%           | 0.179%        |
| 3         | 4.771       | 347           | 354         | 363          | rBV      | 141927         | 202266        | 10.37%          | 1.850%        |
| 4         | 5.258       | 430           | 437         | 446          | rBV      | 589787         | 838854        | 43.02%          | 7.674%        |
| 5         | 6.862       | 704           | 710         | 724          | rBV      | 696571         | 1063425       | 54.54%          | 9.729%        |
| 6         | 7.203       | 761           | 768         | 777          | rBV      | 663141         | 998192        | 51.19%          | 9.132%        |
| 7         | 7.656       | 839           | 845         | 852          | rBV      | 112187         | 174593        | 8.95%           | 1.597%        |
| 8         | 7.979       | 891           | 900         | 906          | rBV      | 378688         | 577864        | 29.64%          | 5.286%        |
| 9         | 8.819       | 1036          | 1043        | 1052         | rBV      | 401797         | 651525        | 33.41%          | 5.960%        |
| 10        | 10.452      | 1313          | 1321        | 1328         | rBV      | 141584         | 231697        | 11.88%          | 2.120%        |
| 11        | 12.932      | 1735          | 1743        | 1750         | rBV      | 752670         | 1143611       | 58.65%          | 10.462%       |
| 12        | 13.778      | 1881          | 1887        | 1895         | rBV      | 147087         | 199546        | 10.23%          | 1.826%        |
| 13        | 14.318      | 1973          | 1979        | 1989         | rVB2     | 203826         | 305292        | 15.66%          | 2.793%        |
| 14        | 15.811      | 2227          | 2233        | 2243         | rBV      | 782647         | 1084340       | 55.61%          | 9.920%        |
| 15        | 17.062      | 2440          | 2446        | 2453         | rBV2     | 284796         | 405760        | 20.81%          | 3.712%        |
| 16        | 17.967      | 2596          | 2600        | 2608         | rBV3     | 23199          | 37714         | 1.93%           | 0.345%        |
| 17        | 19.700      | 2889          | 2895        | 2900         | rBV      | 1637034        | 1949899       | 100.00%         | 17.838%       |
| 18        | 20.963      | 3105          | 3110        | 3117         | rBV2     | 50867          | 76842         | 3.94%           | 0.703%        |
| 19        | 21.251      | 3152          | 3159        | 3163         | rBV      | 376858         | 494581        | 25.36%          | 4.525%        |
| 20        | 23.490      | 3533          | 3540        | 3546         | rBV2     | 239176         | 449489        | 23.05%          | 4.112%        |

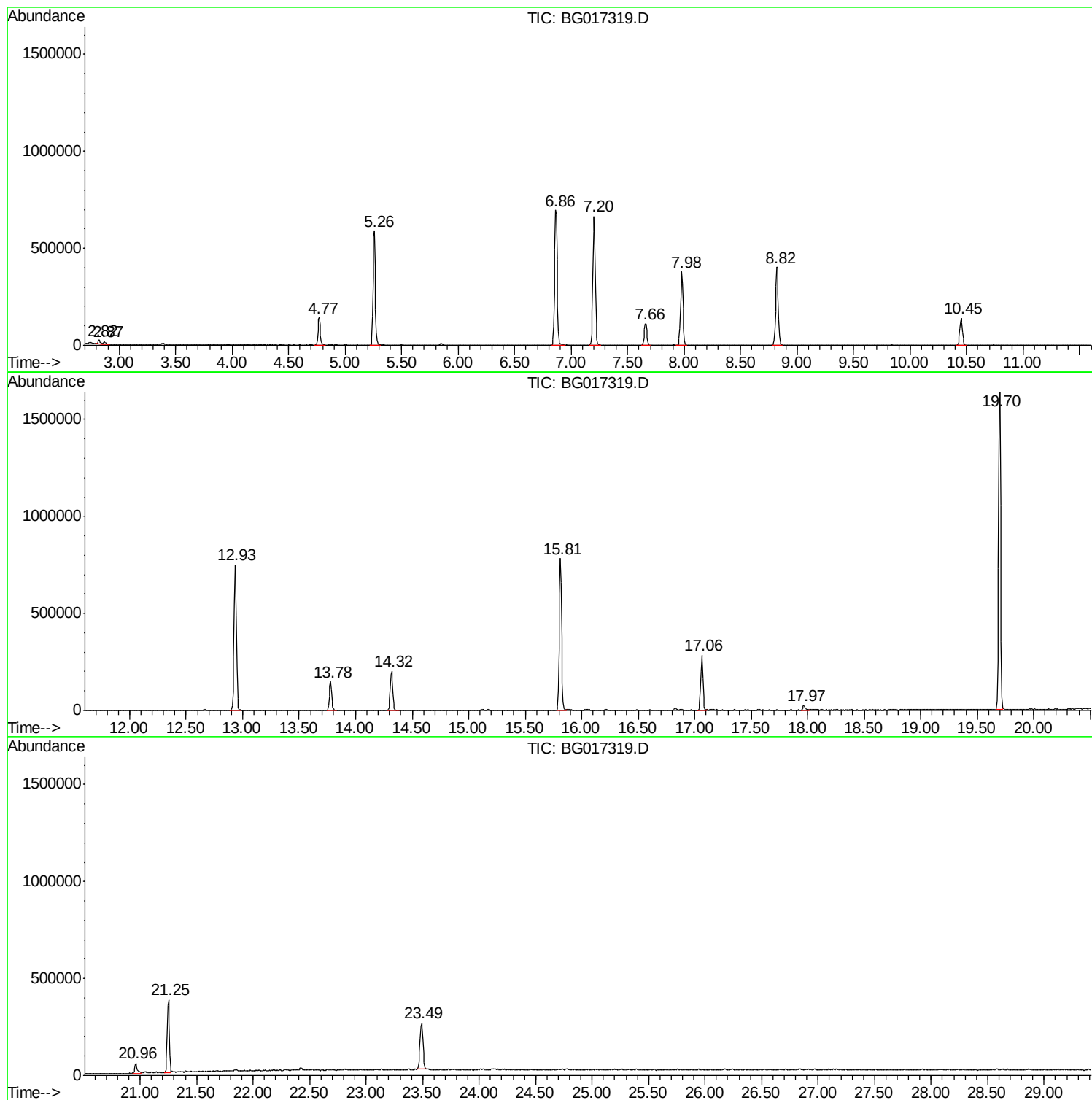
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Operator : TP/IZ  
Sample : G2387-05  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
TP-16-D8

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG051315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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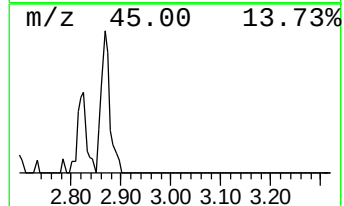
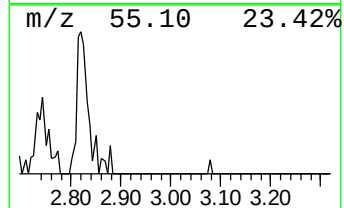
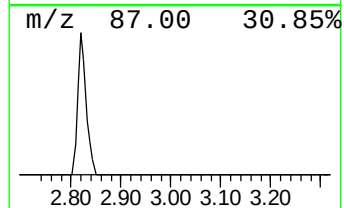
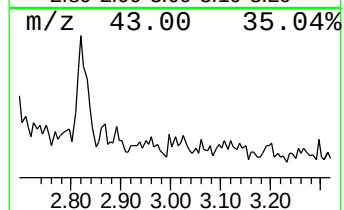
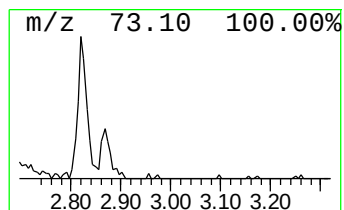
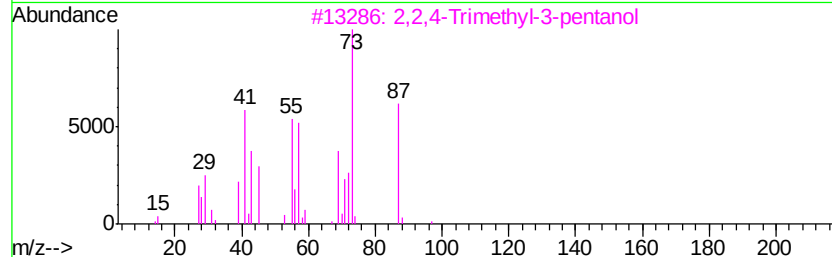
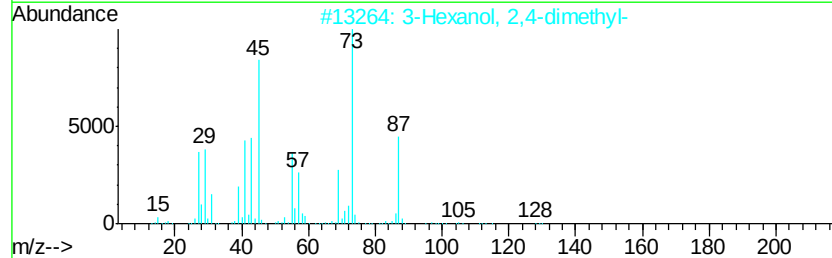
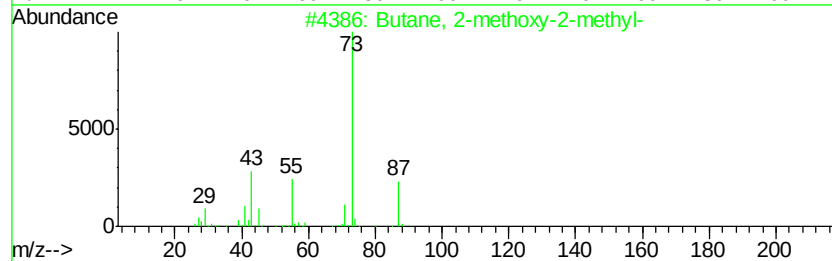
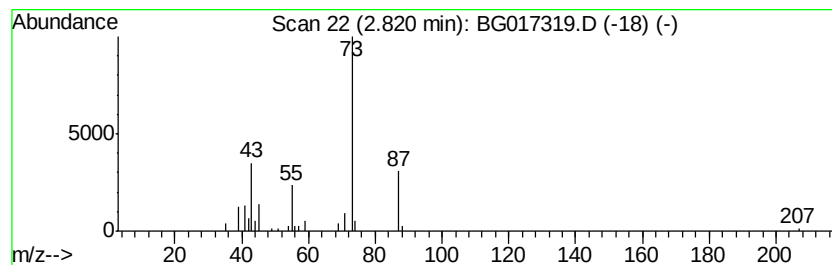
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 2.82 | 2.98 ng | 25990 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl-          | 102 | C6H14O  | 000994-05-8 | 72   |
| 2    |    | 3-Hexanol, 2,4-dimethyl-             | 130 | C8H18O  | 013432-25-2 | 38   |
| 3    |    | 2,2,4-Trimethyl-3-pentanol           | 130 | C8H18O  | 005162-48-1 | 36   |
| 4    |    | Butanoic acid, 2-hydroxy-2-methyl... | 132 | C6H12O3 | 032793-34-3 | 17   |
| 5    |    | Acetamide, N-ethyl-                  | 87  | C4H9NO  | 000625-50-3 | 9    |



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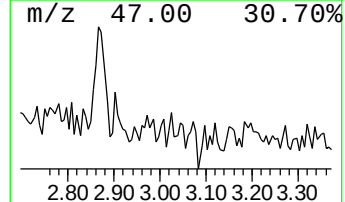
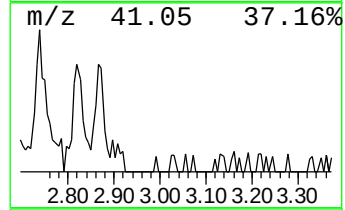
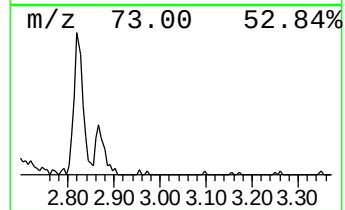
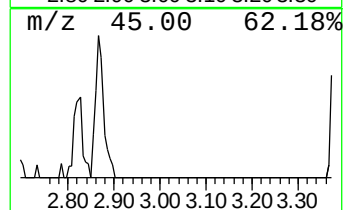
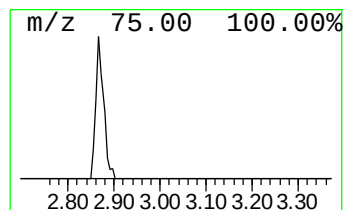
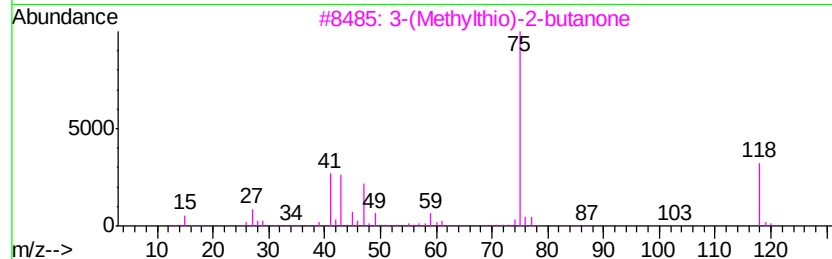
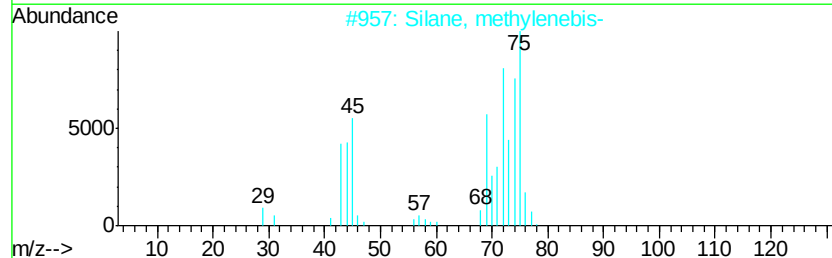
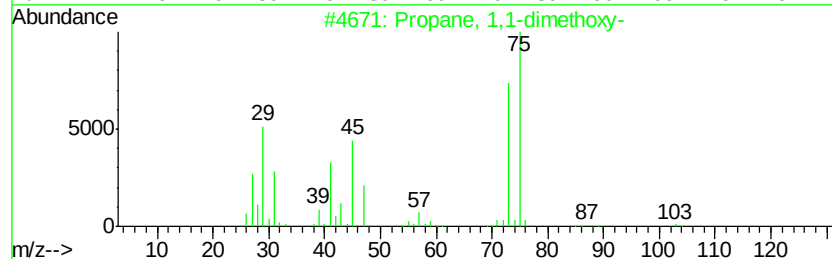
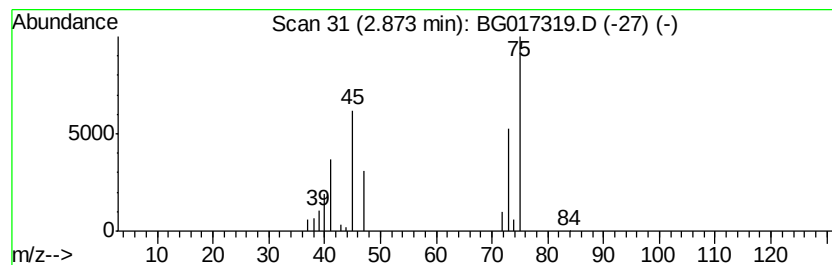
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TIC Integration Parameters: LSCINT.P

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Peak Number 2 Propane, 1,1-dimethoxy- Concentration Rank 6

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 2.87 | 2.24 ng | 19521 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# | of | Tentative ID                            | MW  | MolForm  | CAS#         | Qual |
|------|----|-----------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Propane, 1,1-dimethoxy-                 | 104 | C5H12O2  | 004744-10-9  | 78   |
| 2    |    | Silane, methylenebis-                   | 76  | CH8Si2   | 001759-88-2  | 40   |
| 3    |    | 3-(Methylthio)-2-butanone               | 118 | C5H10OS  | 053475-15-3  | 28   |
| 4    |    | 2-Methylbutane-1,4-diol, 3-(1-ethyl-... | 192 | C9H20O4  | 088481-54-3  | 10   |
| 5    |    | 1-Propanethiol, 3-(diethylborylo...     | 160 | C7H17BO3 | 1000161-78-5 | 9    |



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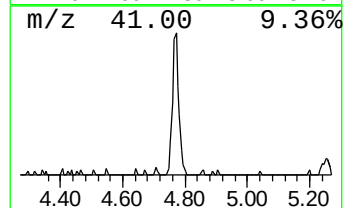
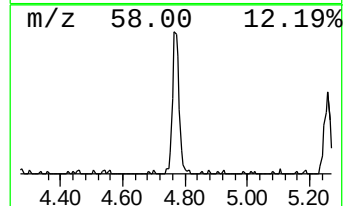
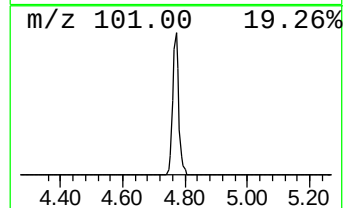
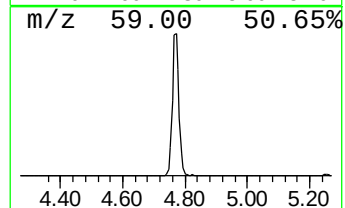
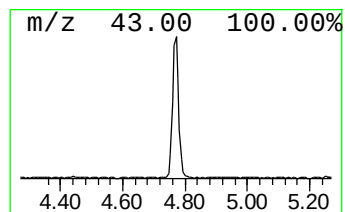
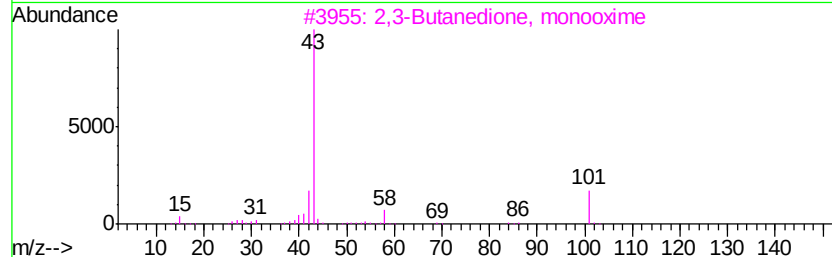
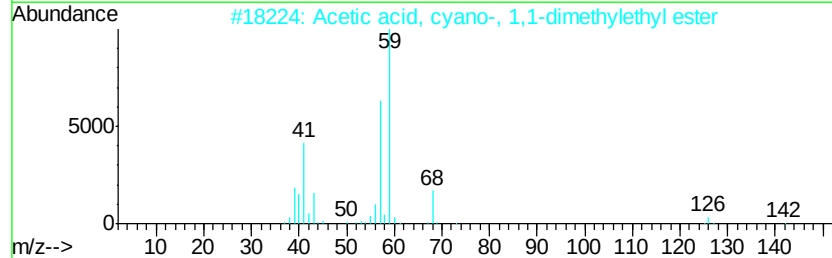
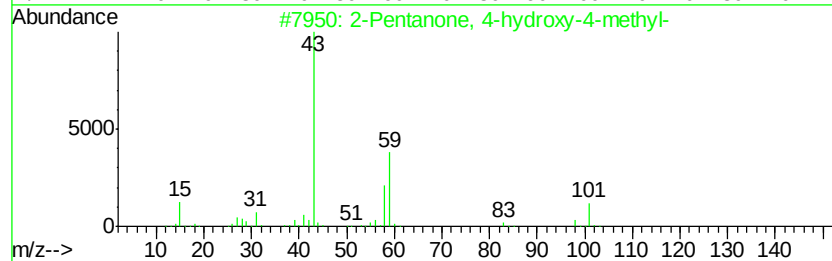
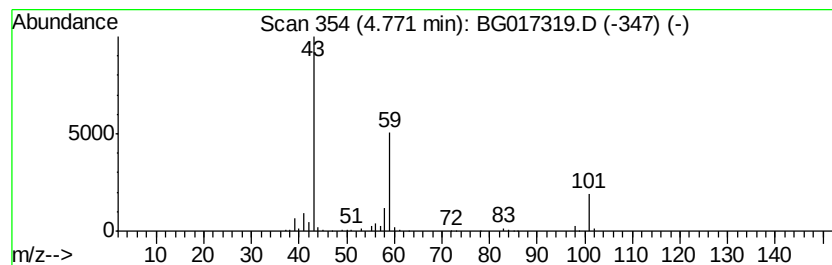
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.77 | 23.17 ng | 202266 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 3    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 9    |
| 4    |    | Morpholine, 4-methyl-                | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    | 3-Hydroxy-3-methyl-2-butanone        | 102 | C5H10O2  | 000115-22-0 | 9    |



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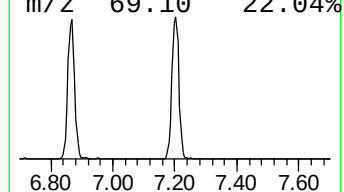
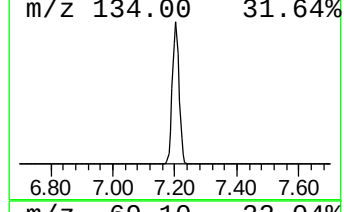
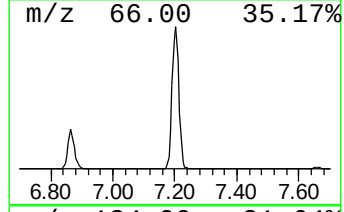
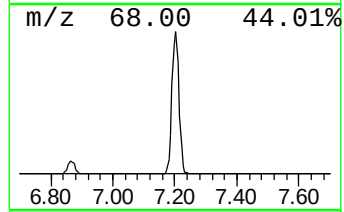
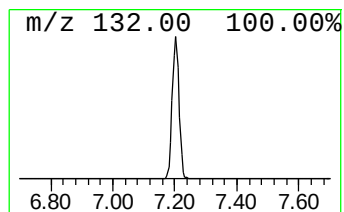
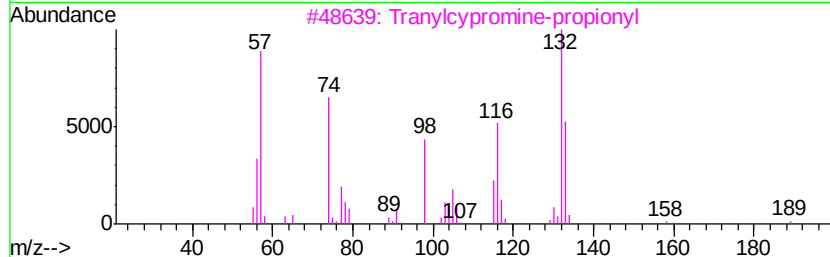
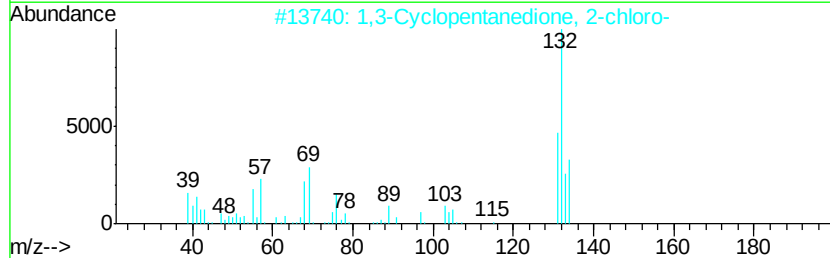
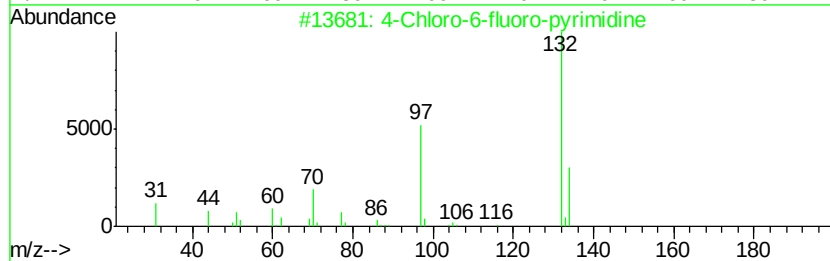
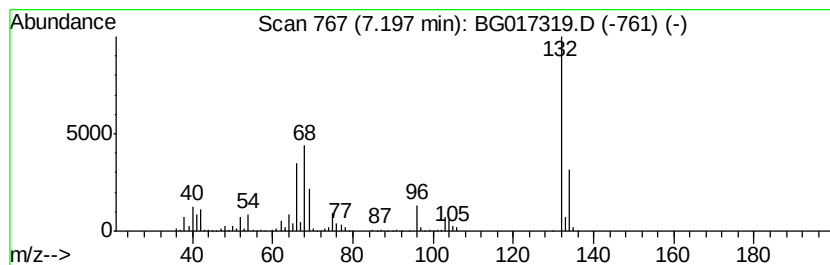
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Peak Number 4 unknown7.20 Concentration Rank 1

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 7.20 | 114.35 ng | 998192 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# of | 5 | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|---------|---|----------------------------------|-----|-----------|--------------|------|
| 1       |   | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2       |   | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 25   |
| 3       |   | Tranvlcyproline-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 22   |
| 4       |   | (5-METHYL-2-PYRIDYL)ACETONITRILE | 132 | C8H8N2    | 1000241-93-9 | 22   |
| 5       |   | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |



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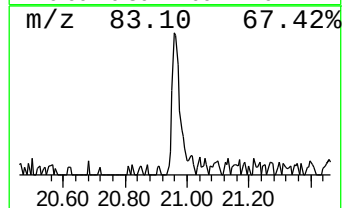
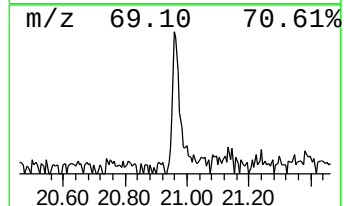
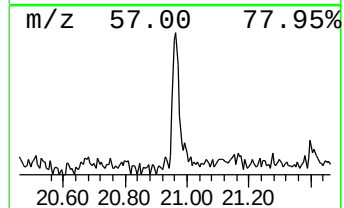
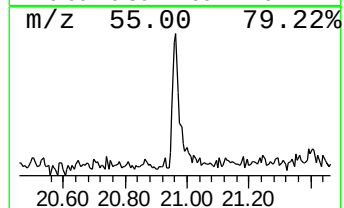
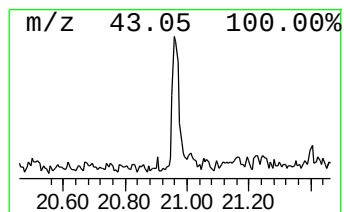
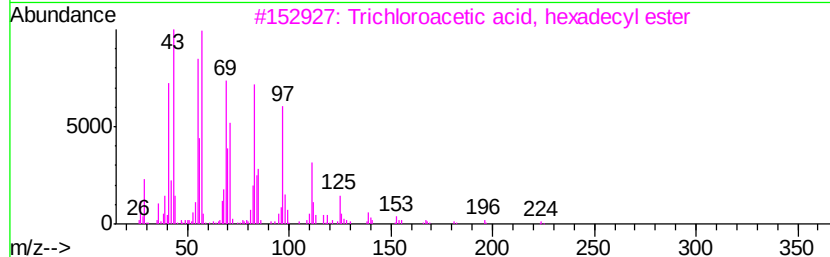
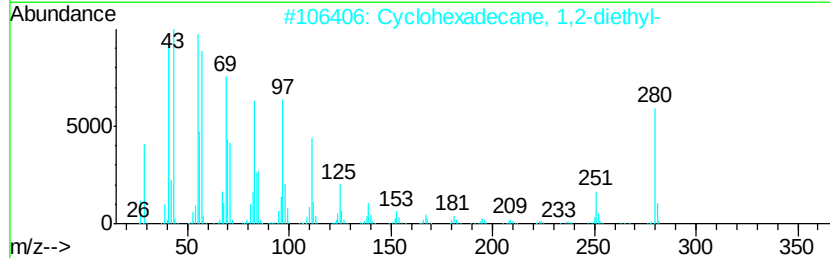
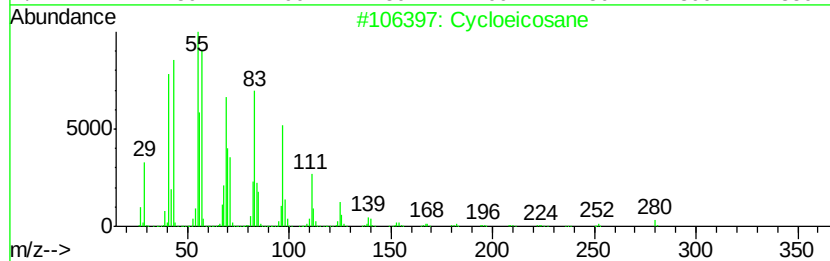
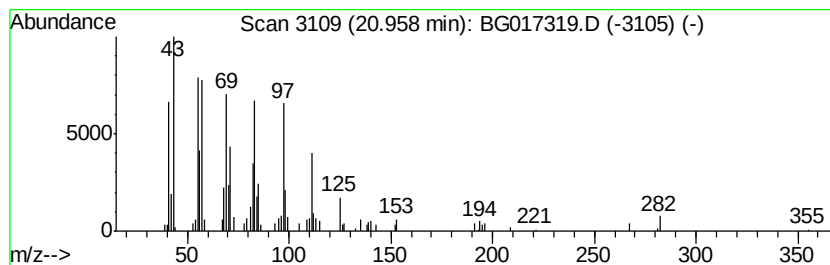
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Peak Number 6 Cycloeicosane Concentration Rank 4

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 20.96   | 3.11 ng | 76842                               | Chrysene-d12     | 21.25           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | Cycloeicosane                       | 280 C20H40       | 000296-56-0 95  |
| 2       |         | Cyclohexadecane, 1,2-diethyl-       | 280 C20H40       | 1000155-85-3 93 |
| 3       |         | Trichloroacetic acid, hexadecyl ... | 386 C18H33Cl3O2  | 074339-54-1 93  |
| 4       |         | Heptafluorobutanoic acid, heptad... | 452 C21H35F7O2   | 1000282-97-3 90 |
| 5       |         | 3-Hexadecene, (Z)-                  | 224 C16H32       | 034303-81-6 90  |



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TP-16-D8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG051315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Butane, 2-methoxy... | 2.82  | 3.0     | ng    | 25990    | 1                     | 7.66  | 174593 | 20.0 |
| Propane, 1,1-dime... | 2.87  | 2.2     | ng    | 19521    | 1                     | 7.66  | 174593 | 20.0 |
| 2-Pentanone, 4-hy... | 4.77  | 23.2    | ng    | 202266   | 1                     | 7.66  | 174593 | 20.0 |
| unknown7.20          | 7.20  | 114.3   | ng    | 998192   | 1                     | 7.66  | 174593 | 20.0 |
| Cycloeicosane        | 20.96 | 3.1     | ng    | 76842    | 5                     | 21.25 | 494581 | 20.0 |