

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA\_F\DATA\BF111117\  
 Data File : BF100456.D  
 Acq On : 11 Nov 2017 23:31  
 Operator : SJ/JU  
 Sample : I6369-01  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 110617-TIDO-F-D-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\_F\METHODS\8270-BF110817.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         | 2.210       | 17            | 21          | 31           | rVB      | 76952          | 103252        | 2.65%           | 0.421%        |
| 2         | 5.122       | 512           | 516         | 525          | rBV      | 289213         | 296892        | 7.62%           | 1.211%        |
| 3         | 5.516       | 579           | 583         | 587          | rBV      | 1625121        | 1477021       | 37.89%          | 6.023%        |
| 4         | 6.516       | 749           | 753         | 757          | rBV      | 1087617        | 1023137       | 26.25%          | 4.172%        |
| 5         | 6.669       | 775           | 779         | 782          | rBV      | 2529373        | 2337535       | 59.97%          | 9.532%        |
| 6         | 6.904       | 815           | 819         | 822          | rBV      | 1022310        | 873277        | 22.40%          | 3.561%        |
| 7         | 7.063       | 841           | 846         | 849          | rBV      | 3077674        | 2945338       | 75.56%          | 12.010%       |
| 8         | 7.469       | 909           | 915         | 918          | rBV      | 2222327        | 2437501       | 62.53%          | 9.939%        |
| 9         | 8.186       | 1032          | 1037        | 1039         | rBV      | 1155469        | 1069589       | 27.44%          | 4.361%        |
| 10        | 8.733       | 1126          | 1130        | 1134         | rBV      | 90325          | 77324         | 1.98%           | 0.315%        |
| 11        | 9.263       | 1214          | 1220        | 1223         | rBV      | 3918439        | 3897824       | 100.00%         | 15.894%       |
| 12        | 9.645       | 1282          | 1285        | 1294         | rBV      | 137002         | 127355        | 3.27%           | 0.519%        |
| 13        | 9.939       | 1329          | 1335        | 1339         | rBV      | 1142925        | 1045995       | 26.84%          | 4.265%        |
| 14        | 10.604      | 1445          | 1448        | 1452         | rVB      | 96878          | 71638         | 1.84%           | 0.292%        |
| 15        | 10.651      | 1452          | 1456        | 1459         | rBV      | 238256         | 207508        | 5.32%           | 0.846%        |
| 16        | 10.727      | 1465          | 1469        | 1472         | rBV      | 1341792        | 1246228       | 31.97%          | 5.082%        |
| 17        | 11.427      | 1584          | 1588        | 1591         | rBV      | 934380         | 846394        | 21.71%          | 3.451%        |
| 18        | 13.010      | 1853          | 1857        | 1861         | rBV      | 2974579        | 2913131       | 74.74%          | 11.879%       |
| 19        | 13.898      | 2004          | 2008        | 2015         | rBV      | 47407          | 62532         | 1.60%           | 0.255%        |
| 20        | 14.063      | 2032          | 2036        | 2045         | rBV      | 916714         | 827929        | 21.24%          | 3.376%        |
| 21        | 15.551      | 2284          | 2289        | 2295         | rBV      | 430992         | 590155        | 15.14%          | 2.406%        |
| 22        | 16.274      | 2409          | 2412        | 2418         | rVB6     | 27730          | 46004         | 1.18%           | 0.188%        |

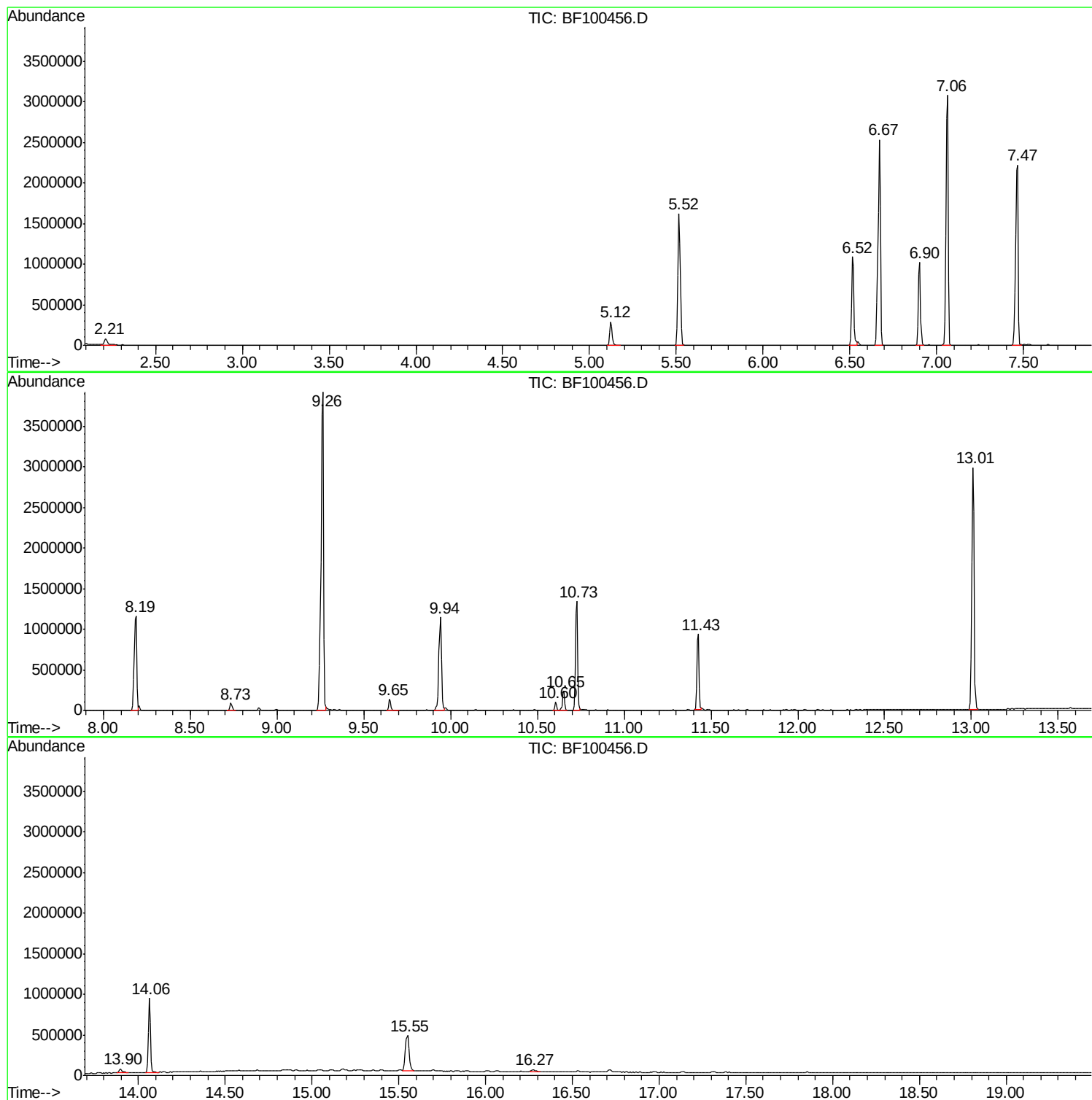
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ClientSampleId :  
110617-TIDO-F-D-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF110817.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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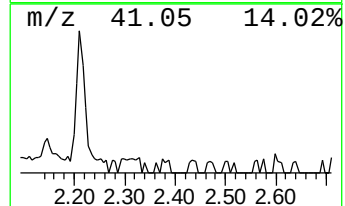
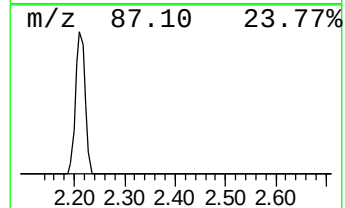
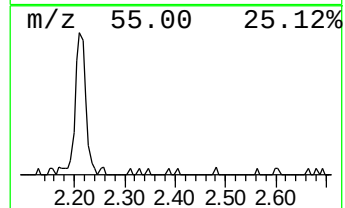
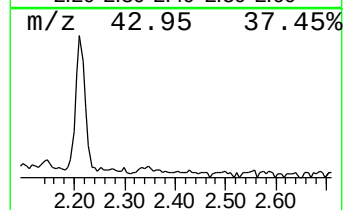
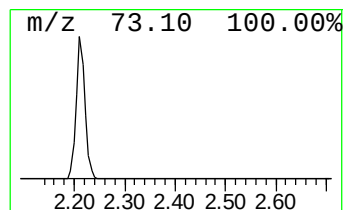
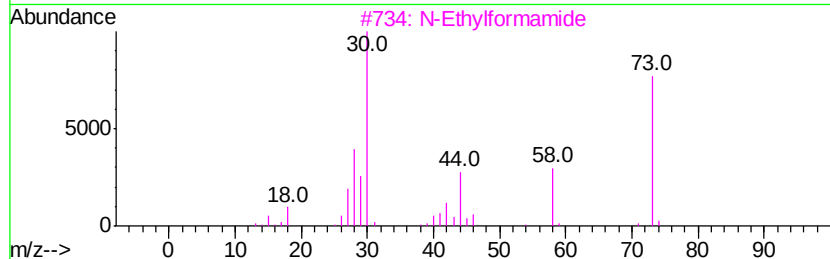
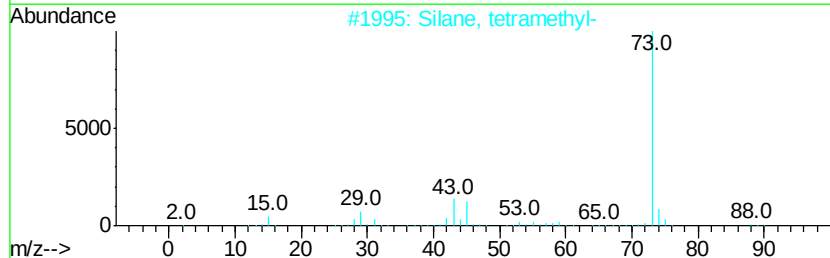
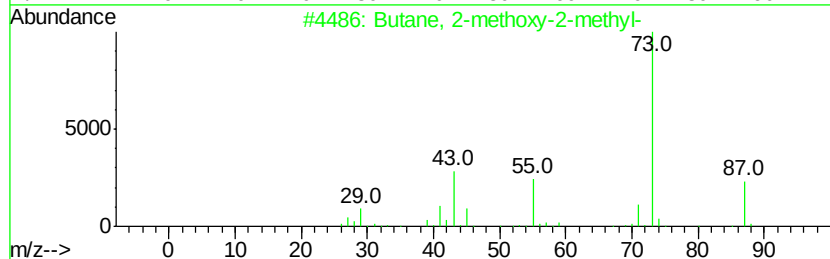
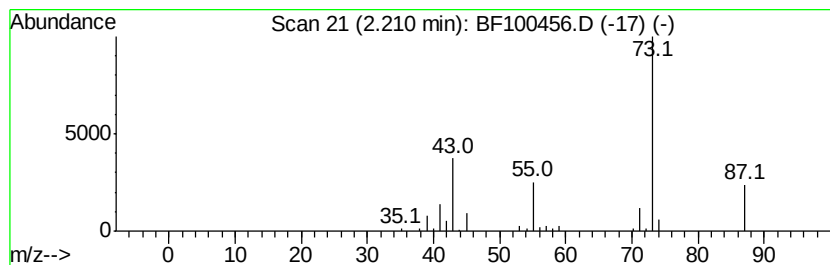
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 2.21 | 2.36 ng | 103252 | 1,4-Dichlorobenzene-d4 | 6.90 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 25   |
| 3    |    |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 4    |    |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |
| 5    |    |   | 3-Hexanol, 2-methyl-        | 116 | C7H16O  | 000617-29-8 | 9    |



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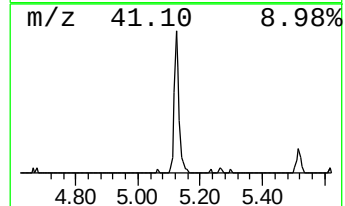
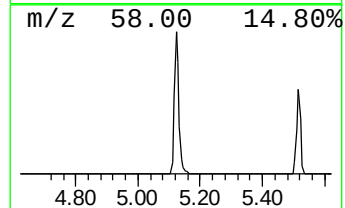
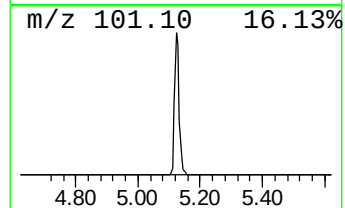
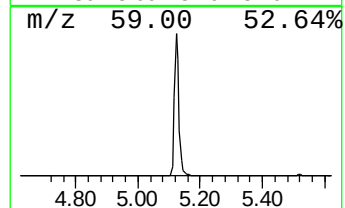
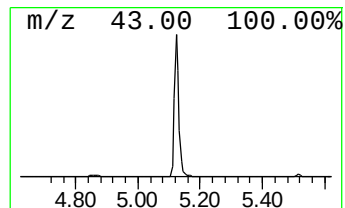
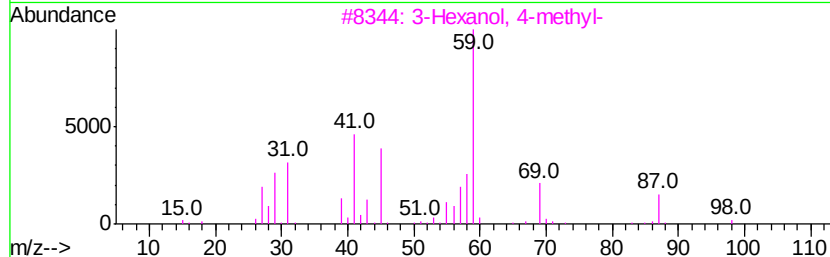
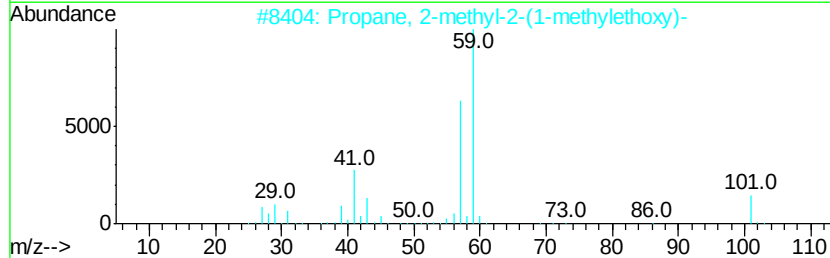
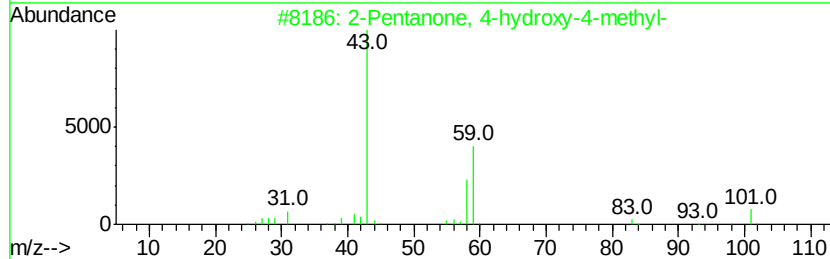
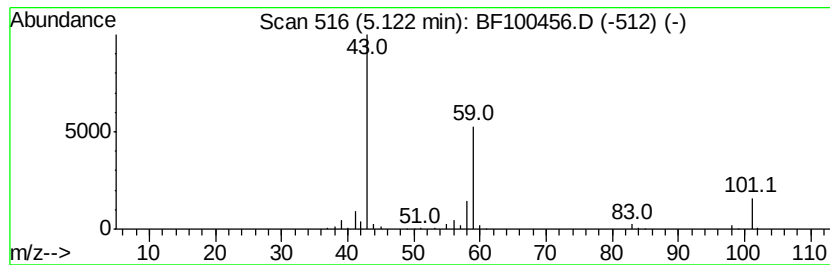
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.12 | 6.80 ng | 296892 | 1,4-Dichlorobenzene-d4 | 6.90 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O  | 017348-59-3 | 42   |
| 3    |    | 3-Hexanol, 4-methyl-                | 116 | C7H16O  | 000615-29-2 | 33   |
| 4    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 28   |
| 5    |    | 2-Hexanol, 2-methyl-                | 116 | C7H16O  | 000625-23-0 | 28   |



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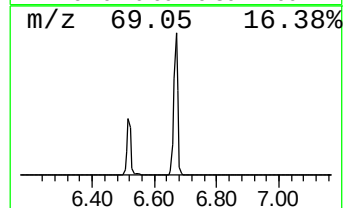
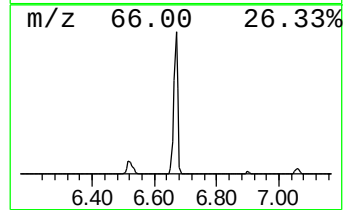
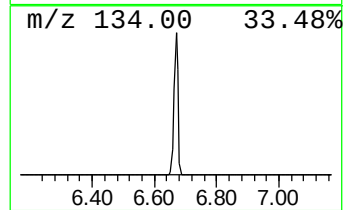
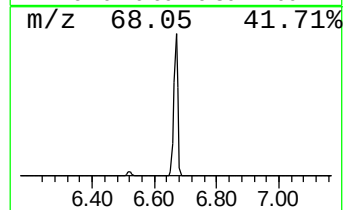
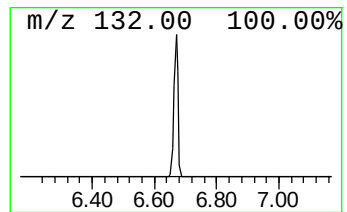
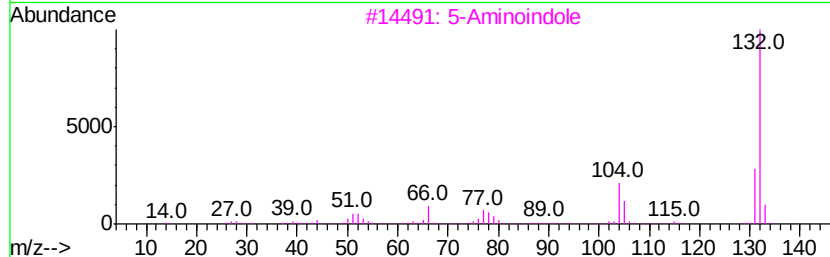
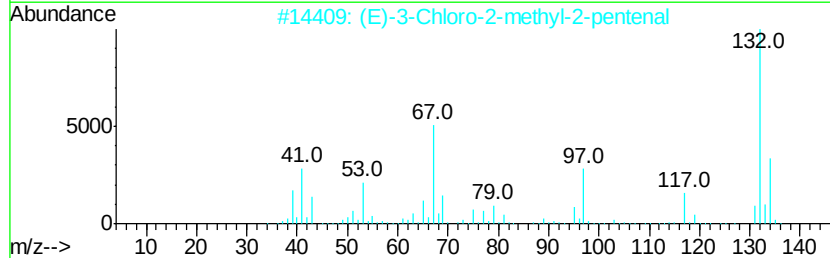
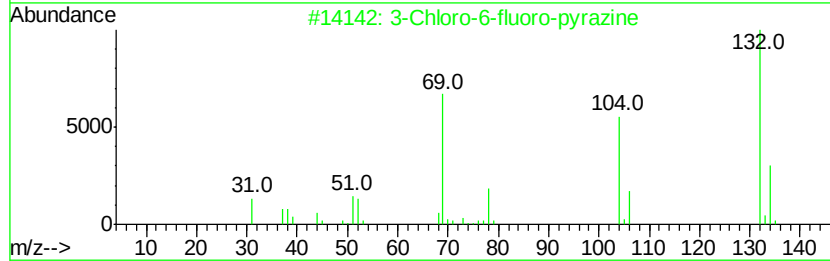
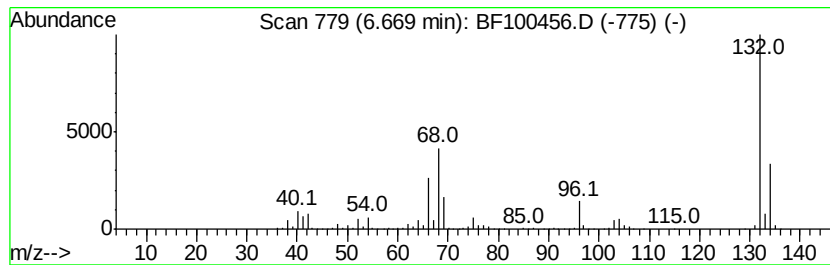
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.67 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.67 | 53.53 ng | 2337540 | 1,4-Dichlorobenzene-d4 | 6.90 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 3    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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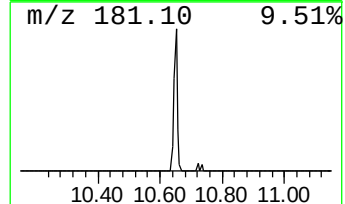
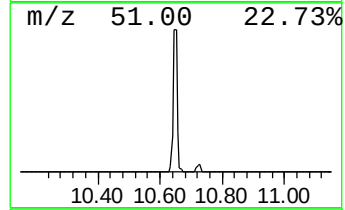
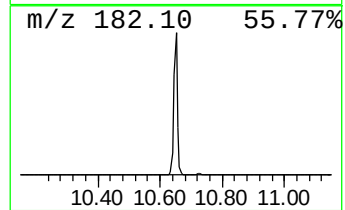
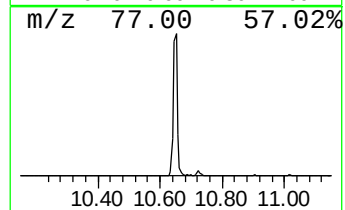
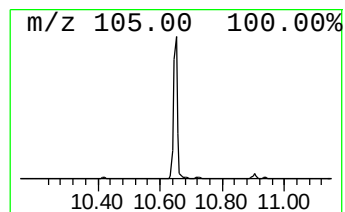
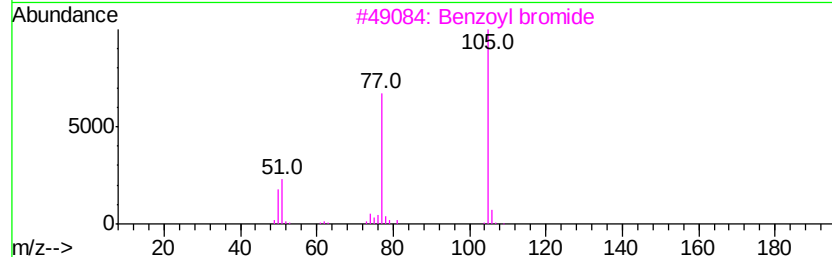
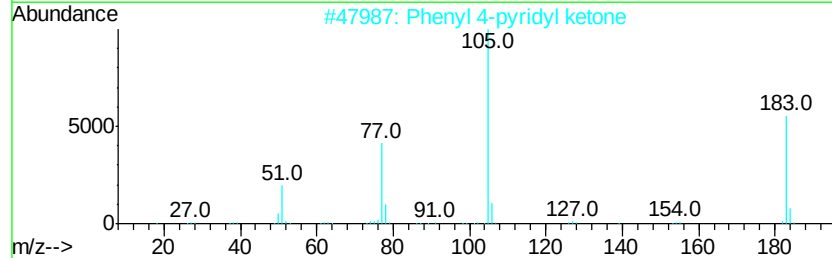
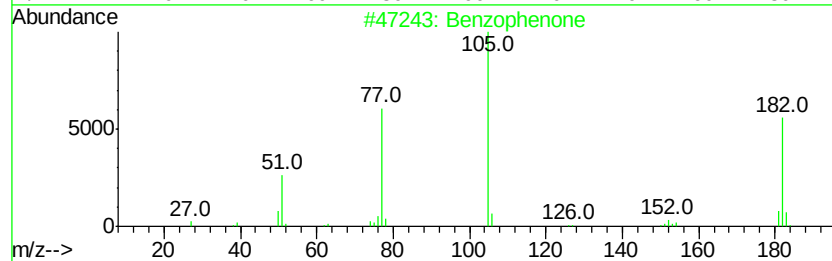
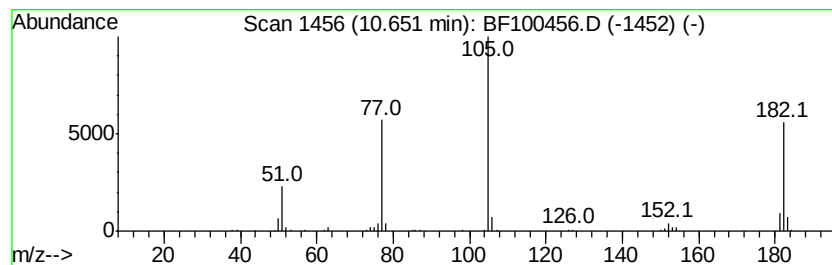
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Peak Number 5 Benzophenone Concentration Rank 4

| R.T.    | EstConc | Area                               | Relative to ISTD | R.T.           |
|---------|---------|------------------------------------|------------------|----------------|
| 10.65   | 3.97 ng | 207508                             | Acenaphthene-d10 | 9.94           |
| Hit# of | 5       | Tentative ID                       | MW MolForm       | CAS# Qual      |
| 1       |         | Benzophenone                       | 182 C13H10O      | 000119-61-9 96 |
| 2       |         | Phenyl 4-pyridyl ketone            | 183 C12H9NO      | 014548-46-0 47 |
| 3       |         | Benzoyl bromide                    | 184 C7H5BrO      | 000618-32-6 47 |
| 4       |         | Vinyl benzoate                     | 148 C9H8O2       | 000769-78-8 47 |
| 5       |         | Benzenebutanoic acid, .gamma.-oxo- | 178 C10H10O3     | 002051-95-8 47 |



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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |      |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|------|---------|------|
|                      |       |         |       |          | #                     | RT   | Resp    | Conc |
| Butane, 2-methoxy... | 2.21  | 2.4     | ng    | 103252   | 1                     | 6.90 | 873277  | 20.0 |
| 2-Pentanone, 4-hy... | 5.12  | 6.8     | ng    | 296892   | 1                     | 6.90 | 873277  | 20.0 |
| unknown6.67          | 6.67  | 53.5    | ng    | 2337540  | 1                     | 6.90 | 873277  | 20.0 |
| Benzophenone         | 10.65 | 4.0     | ng    | 207508   | 3                     | 9.94 | 1046000 | 20.0 |