

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF042116\  
 Data File : BF086348.D  
 Acq On : 21 Apr 2016 17:22  
 Operator : UM/SJ  
 Sample : PB89669BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB89669BL

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-BF042016.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         | 2.190       | 7             | 9           | 15           | rVB2     | 47960          | 100508        | 1.49%           | 0.171%        |
| 2         | 4.407       | 200           | 203         | 211          | rBV      | 244760         | 415565        | 6.16%           | 0.709%        |
| 3         | 5.047       | 257           | 259         | 269          | rBV      | 1264375        | 2224763       | 33.00%          | 3.796%        |
| 4         | 5.459       | 292           | 295         | 298          | rBV      | 3933318        | 5993704       | 88.90%          | 10.227%       |
| 5         | 5.859       | 329           | 330         | 333          | rBV      | 111040         | 107899        | 1.60%           | 0.184%        |
| 6         | 5.939       | 333           | 337         | 345          | rVB2     | 21560          | 82878         | 1.23%           | 0.141%        |
| 7         | 6.487       | 382           | 385         | 394          | rBV      | 4294816        | 5512612       | 81.76%          | 9.406%        |
| 8         | 6.624       | 394           | 397         | 399          | rBV      | 5023900        | 5745979       | 85.22%          | 9.804%        |
| 9         | 6.853       | 415           | 417         | 428          | rBV      | 1460929        | 1565821       | 23.22%          | 2.672%        |
| 10        | 7.013       | 428           | 431         | 433          | rBV      | 4062699        | 4675201       | 69.34%          | 7.977%        |
| 11        | 7.413       | 463           | 466         | 469          | rBV      | 2580121        | 3665949       | 54.37%          | 6.255%        |
| 12        | 8.133       | 526           | 529         | 532          | rBV      | 1433495        | 1969787       | 29.22%          | 3.361%        |
| 13        | 8.259       | 538           | 540         | 551          | rVB      | 34727          | 100231        | 1.49%           | 0.171%        |
| 14        | 9.219       | 620           | 624         | 626          | rBV      | 5319514        | 6742192       | 100.00%         | 11.504%       |
| 15        | 9.893       | 680           | 683         | 685          | rBV      | 2181672        | 2349285       | 34.84%          | 4.009%        |
| 16        | 10.682      | 749           | 752         | 765          | rBV      | 2990174        | 4180510       | 62.01%          | 7.133%        |
| 17        | 11.379      | 810           | 813         | 825          | rBV      | 2048464        | 2695800       | 39.98%          | 4.600%        |
| 18        | 12.968      | 949           | 952         | 954          | rBV      | 5174732        | 6567613       | 97.41%          | 11.206%       |
| 19        | 14.008      | 1040          | 1043        | 1054         | rBV      | 1955148        | 2339794       | 34.70%          | 3.992%        |
| 20        | 15.437      | 1165          | 1168        | 1175         | rBV      | 858887         | 1570282       | 23.29%          | 2.679%        |

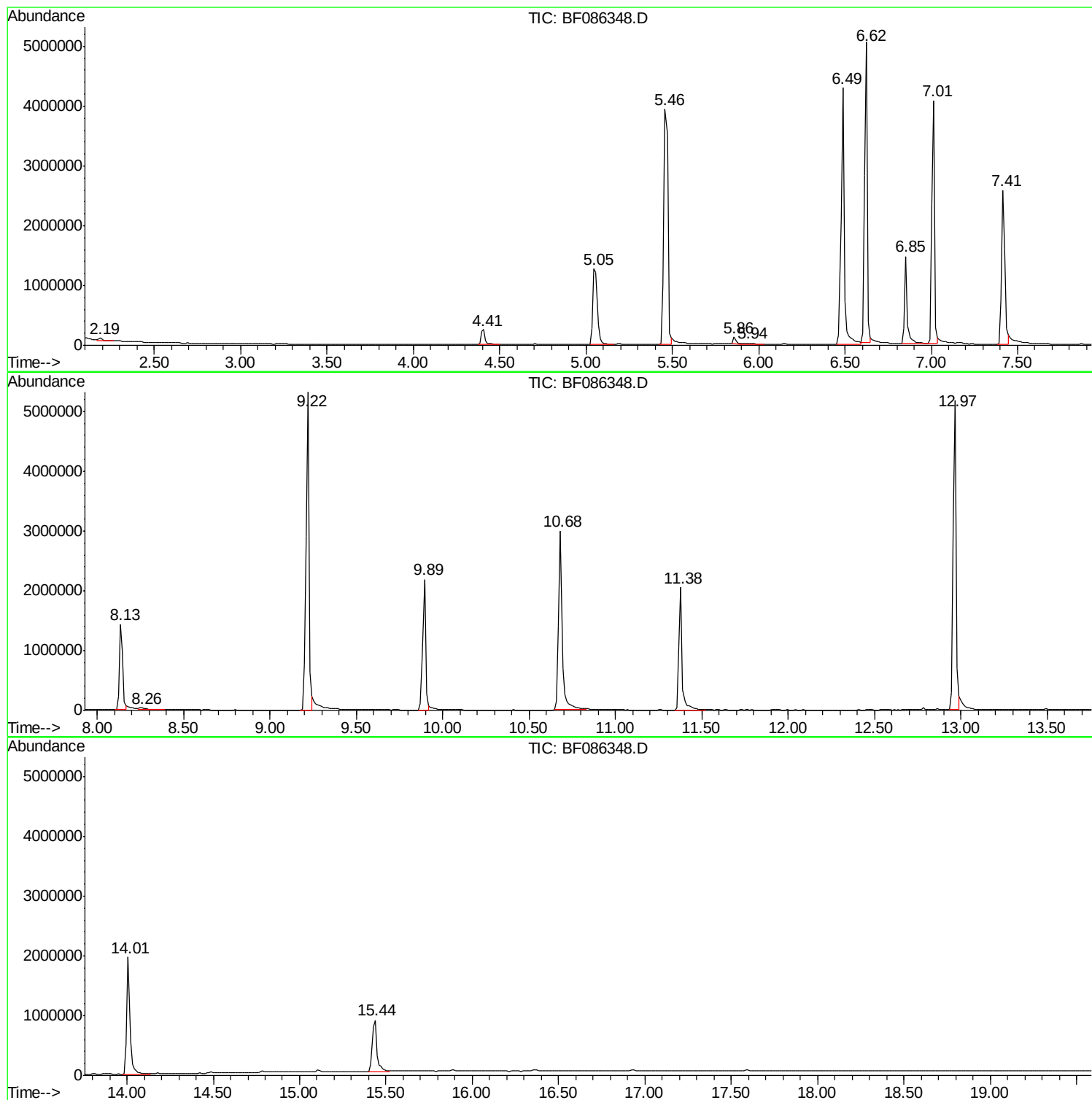
Sum of corrected areas: 58606373

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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF042016.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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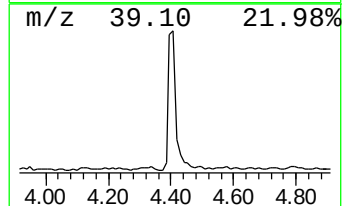
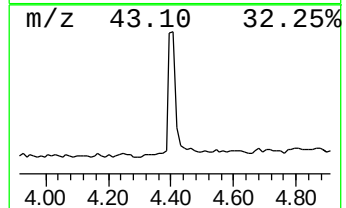
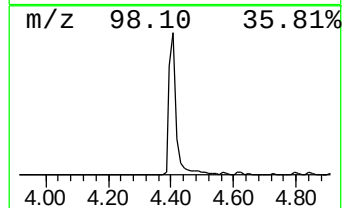
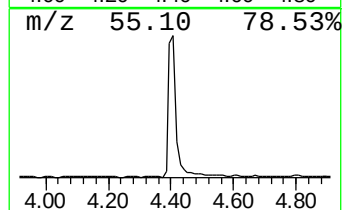
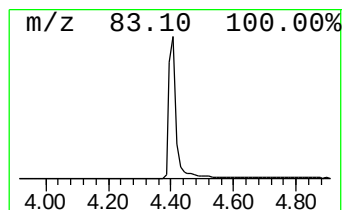
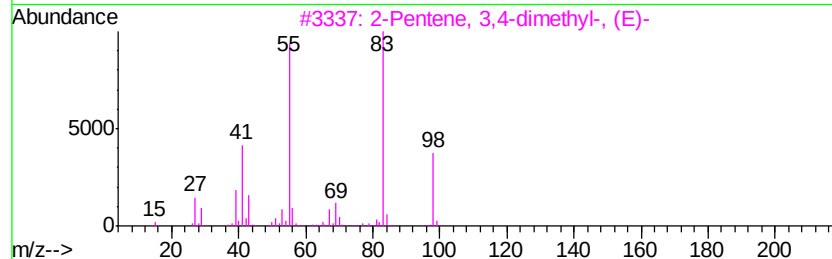
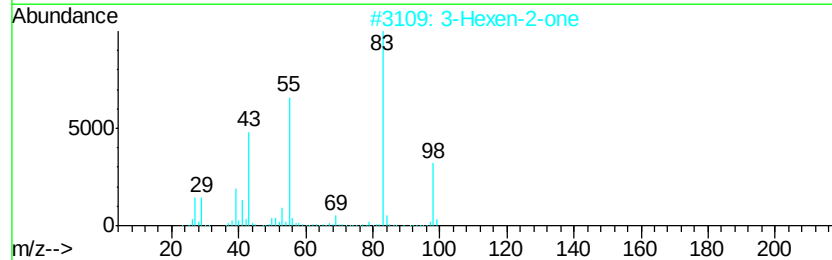
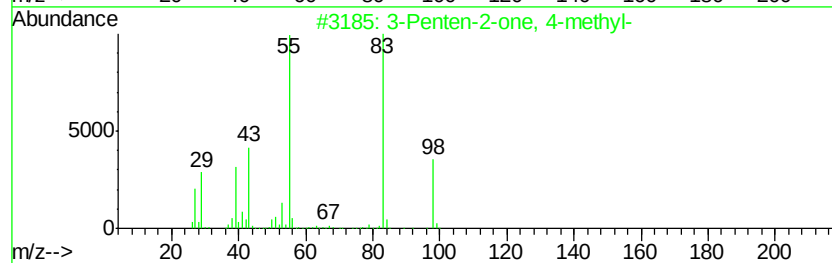
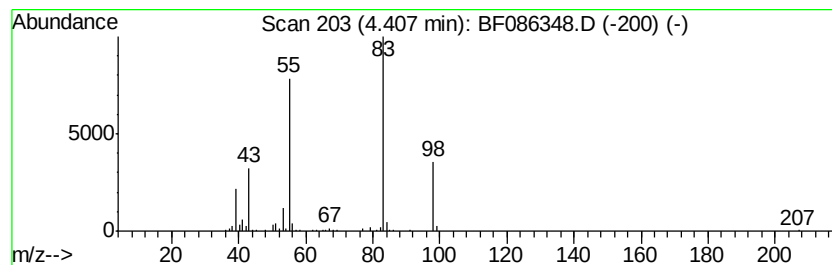
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.41 | 5.31 ng | 415565 | 1,4-Dichlorobenzene-d4 | 6.85 |

| Hit# | of | 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 91   |
| 2    |    |   | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 91   |
| 3    |    |   | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 86   |
| 4    |    |   | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 78   |
| 5    |    |   | Furan, 2-methoxy-              | 98 | C5H6O2  | 025414-22-6 | 78   |



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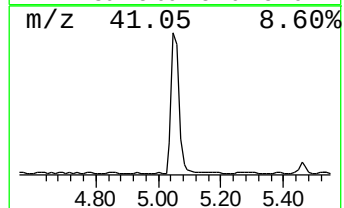
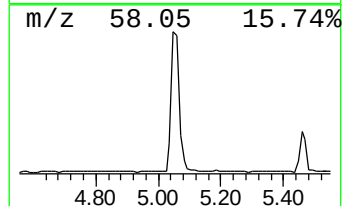
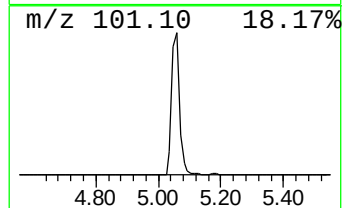
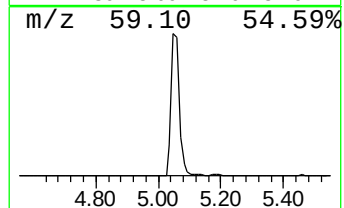
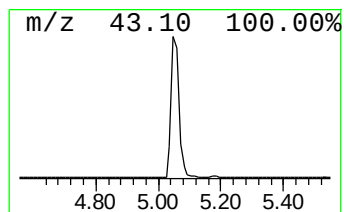
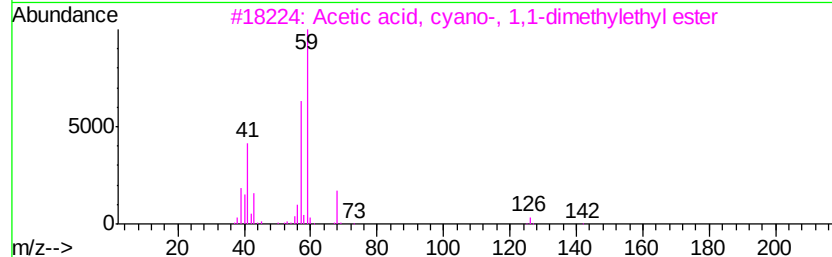
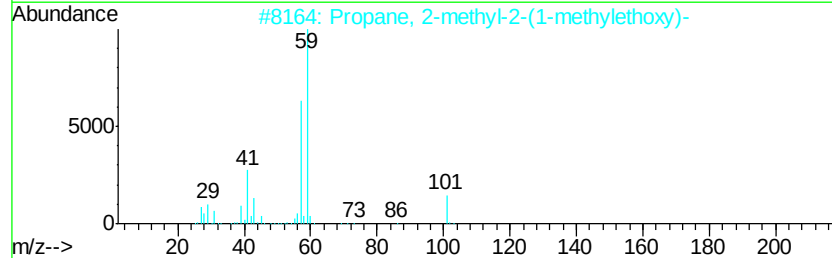
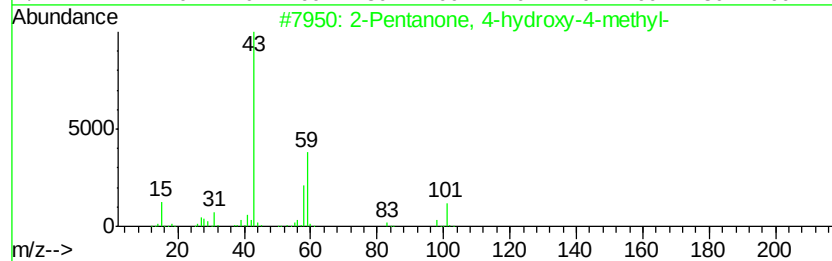
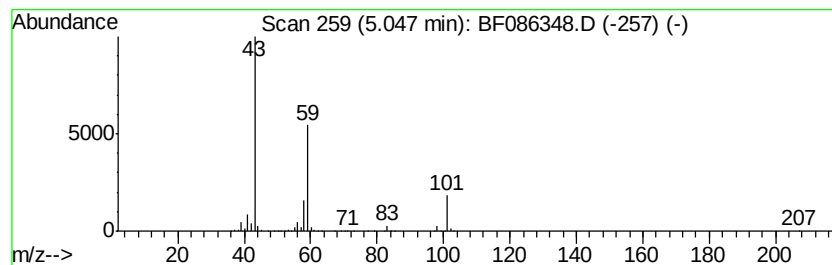
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.05 | 28.42 ng | 2224760 | 1,4-Dichlorobenzene-d4 | 6.85 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    | Propane, 2-methyl-2-(1-methyleth...  | 116 | C7H16O   | 017348-59-3 | 33   |
| 3    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 4    |    | 1-Propen-2-ol, acetate               | 100 | C5H8O2   | 000108-22-5 | 10   |
| 5    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 9    |



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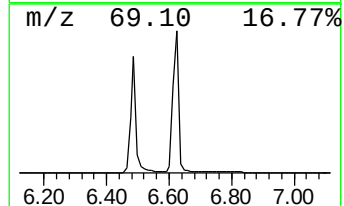
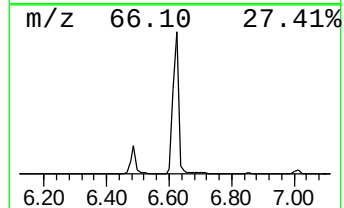
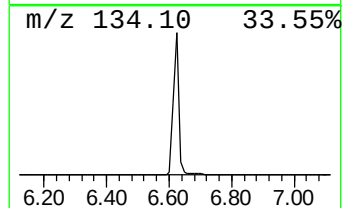
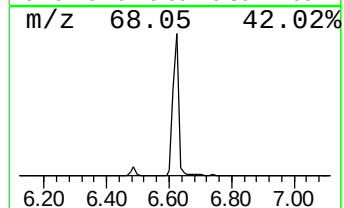
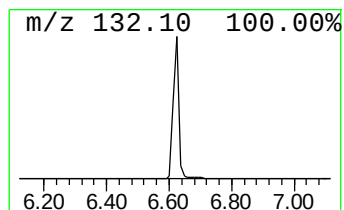
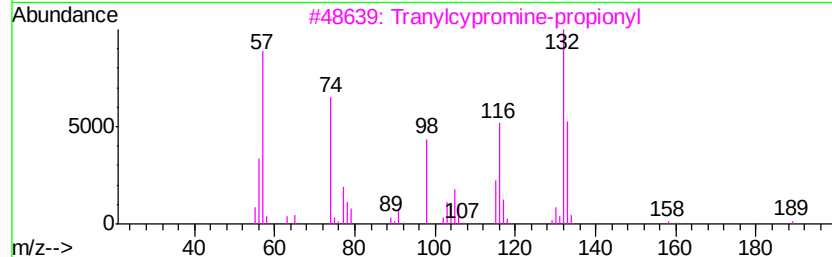
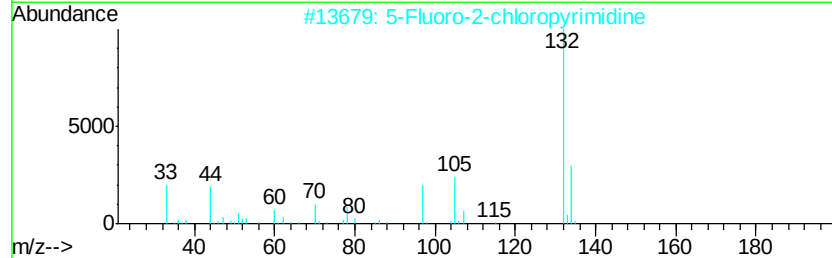
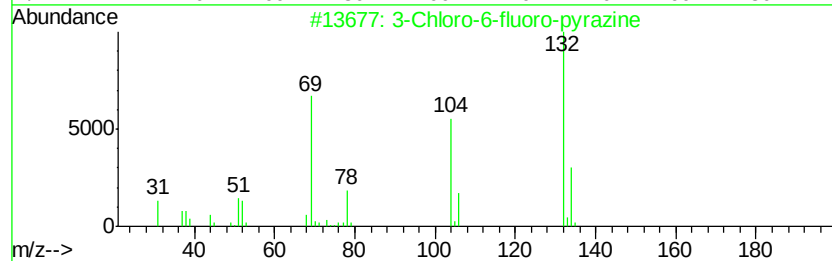
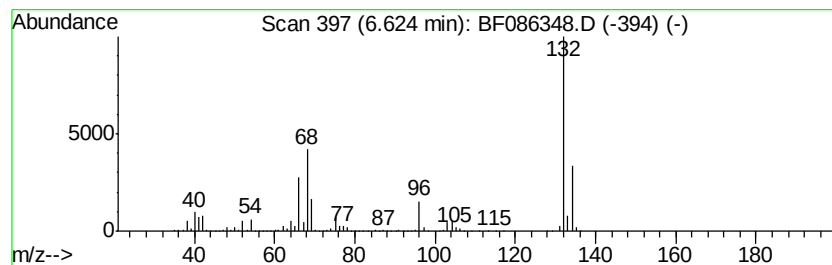
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Peak Number 3 unknown6.62 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.62 | 73.39 ng | 5745980 | 1,4-Dichlorobenzene-d4 | 6.85 |

| Hit# | of | Tentative ID                | MW  | MolForm   | CAS#         | Qual |
|------|----|-----------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine  | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | 5-Fluoro-2-chloropyrimidine | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 3    |    | Tranvlcypropine-propionyl   | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | Benzene, 1,2,4-trifluoro-   | 132 | C6H3F3    | 000367-23-7  | 9    |
| 5    |    | Benzene, 1,3,5-trifluoro-   | 132 | C6H3F3    | 000372-38-3  | 9    |



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| TIC Top Hit name     | RT   | EstConc | Units | Response | --Internal Standard-- |      |         |      |
|----------------------|------|---------|-------|----------|-----------------------|------|---------|------|
|                      |      |         |       |          | #                     | RT   | Resp    | Conc |
| 3-Penten-2-one, 4... | 4.41 | 5.3     | ng    | 415565   | 1                     | 6.85 | 1565820 | 20.0 |
| 2-Pentanone, 4-hy... | 5.05 | 28.4    | ng    | 2224760  | 1                     | 6.85 | 1565820 | 20.0 |
| unknown6.62          | 6.62 | 73.4    | ng    | 5745980  | 1                     | 6.85 | 1565820 | 20.0 |