

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF103017\  
 Data File : BF099986.D  
 Acq On : 31 Oct 2017 3:45  
 Operator : SJ/JU  
 Sample : I6083-12  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 102017-SIDI-F-D-M

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\_F\METHODS\8270-BF102317.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         | 4.999       | 492           | 495         | 512          | rBV      | 46951          | 76808         | 3.29%           | 0.556%        |
| 2         | 5.410       | 562           | 565         | 587          | rBV      | 399564         | 595819        | 25.49%          | 4.316%        |
| 3         | 6.434       | 736           | 739         | 754          | rBV      | 254372         | 407786        | 17.45%          | 2.954%        |
| 4         | 6.557       | 757           | 760         | 773          | rVB      | 1297695        | 1187336       | 50.80%          | 8.602%        |
| 5         | 6.787       | 796           | 799         | 809          | rBV      | 466152         | 428917        | 18.35%          | 3.107%        |
| 6         | 6.940       | 822           | 825         | 847          | rVB      | 1791276        | 1629062       | 69.69%          | 11.802%       |
| 7         | 7.357       | 892           | 896         | 915          | rBV      | 1266365        | 1203750       | 51.50%          | 8.720%        |
| 8         | 8.069       | 1014          | 1017        | 1035         | rBV      | 694097         | 587770        | 25.15%          | 4.258%        |
| 9         | 8.634       | 1110          | 1113        | 1122         | rBB      | 29659          | 32688         | 1.40%           | 0.237%        |
| 10        | 9.145       | 1195          | 1200        | 1203         | rBV      | 2686052        | 2337480       | 100.00%         | 16.934%       |
| 11        | 9.545       | 1265          | 1268        | 1273         | rBV      | 39469          | 37392         | 1.60%           | 0.271%        |
| 12        | 9.822       | 1312          | 1315        | 1329         | rVB      | 683010         | 659728        | 28.22%          | 4.779%        |
| 13        | 10.492      | 1427          | 1429        | 1434         | rBV      | 32632          | 26911         | 1.15%           | 0.195%        |
| 14        | 10.539      | 1434          | 1437        | 1444         | rBV      | 159474         | 127722        | 5.46%           | 0.925%        |
| 15        | 10.622      | 1447          | 1451        | 1465         | rBV      | 886192         | 850092        | 36.37%          | 6.158%        |
| 16        | 11.316      | 1566          | 1569        | 1579         | rBV      | 665590         | 625663        | 26.77%          | 4.533%        |
| 17        | 12.910      | 1835          | 1840        | 1843         | rBV      | 2066630        | 2093162       | 89.55%          | 15.164%       |
| 18        | 13.786      | 1986          | 1989        | 1998         | rBV2     | 28197          | 40577         | 1.74%           | 0.294%        |
| 19        | 13.980      | 2018          | 2022        | 2032         | rBV      | 427764         | 456497        | 19.53%          | 3.307%        |
| 20        | 15.468      | 2271          | 2275        | 2286         | rVB      | 295295         | 398537        | 17.05%          | 2.887%        |

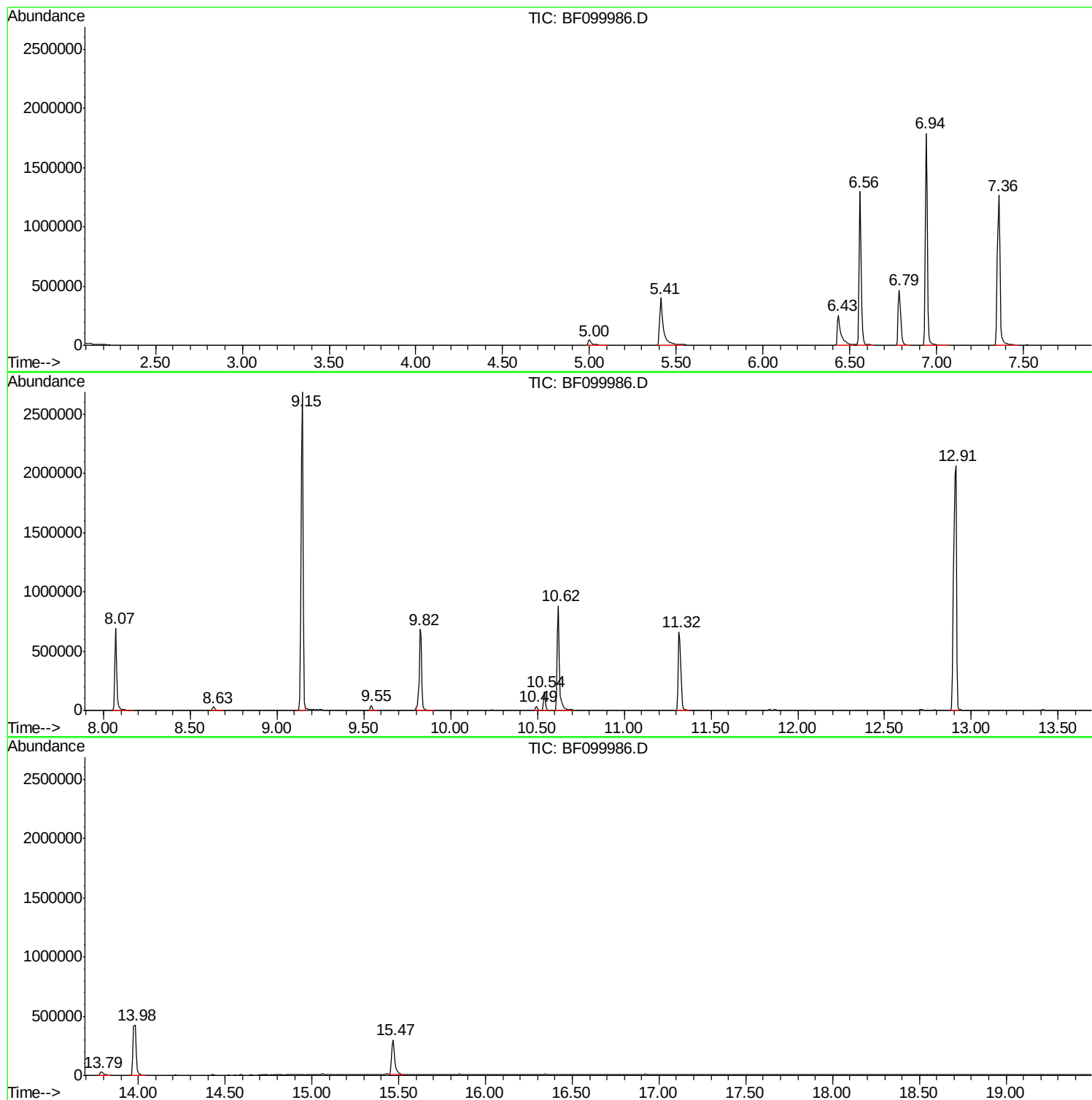
Sum of corrected areas: 13803697

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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF102317.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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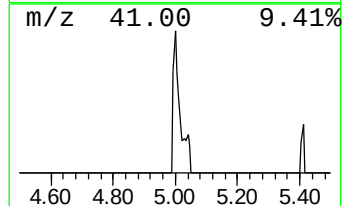
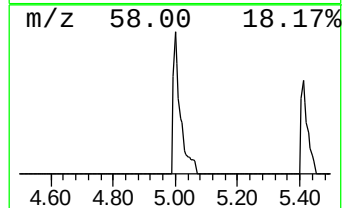
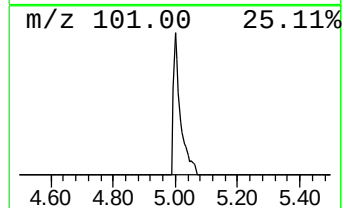
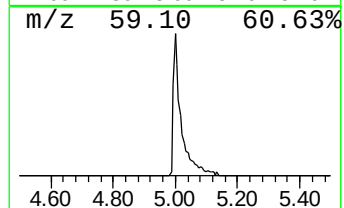
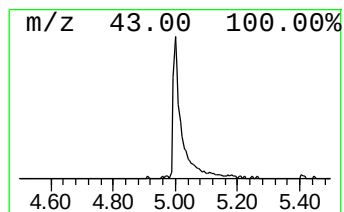
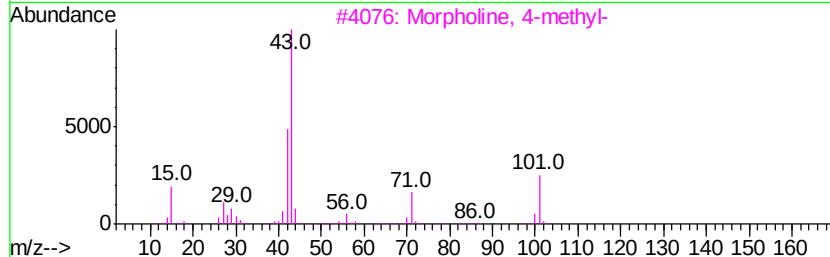
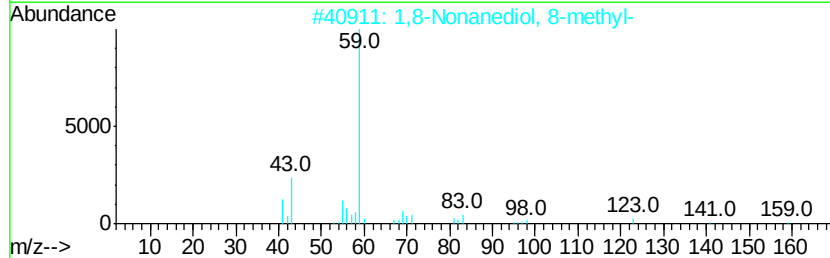
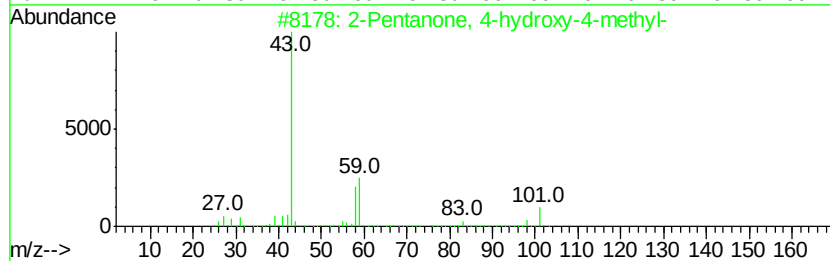
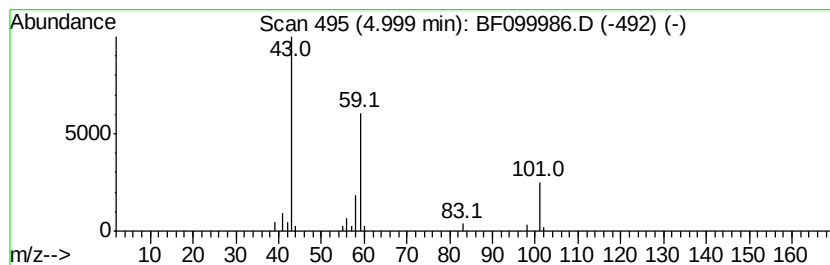
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TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.00 | 3.58 ng | 76808 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    |   | 1,8-Nonanediol, 8-methyl-        | 174 | C10H22O2 | 054725-73-4 | 25   |
| 3    |    |   | Morpholine, 4-methyl-            | 101 | C5H11NO  | 000109-02-4 | 9    |
| 4    |    |   | 2,3-Butanedione, monooxime       | 101 | C4H7NO2  | 000057-71-6 | 9    |
| 5    |    |   | 5-Hexen-2-one                    | 98  | C6H10O   | 000109-49-9 | 9    |



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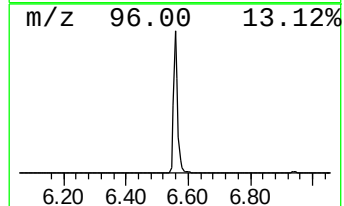
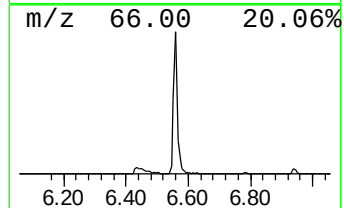
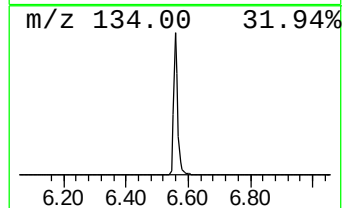
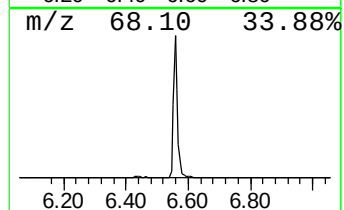
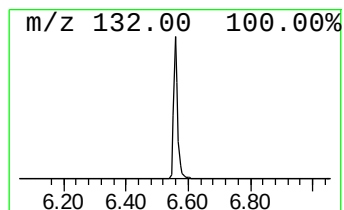
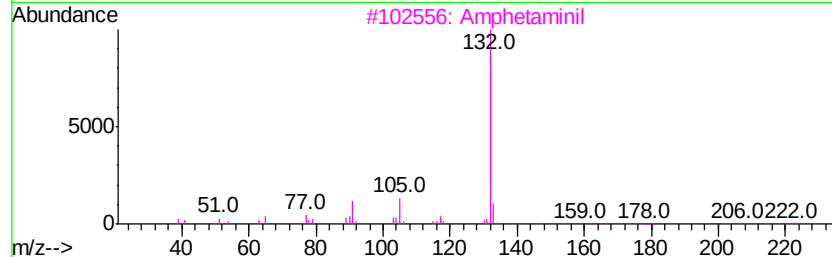
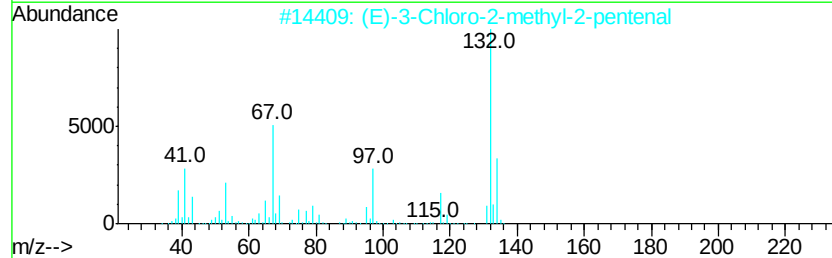
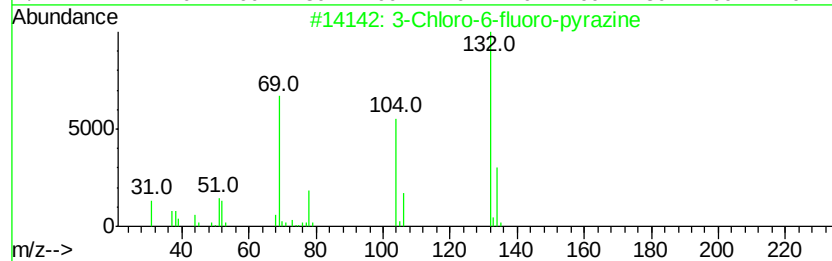
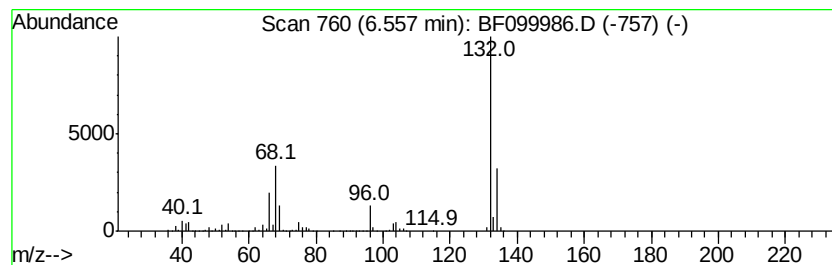
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Peak Number 2 unknown6.56 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.56 | 55.36 ng | 1187340 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 12   |
| 4    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |
| 5    |    | Imidazole, 4,5-dicyano-1-methyl- | 132 | C6H4N4    | 019485-35-9  | 9    |



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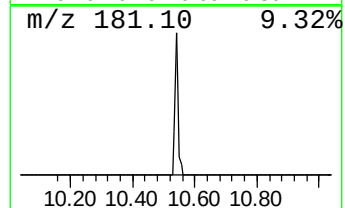
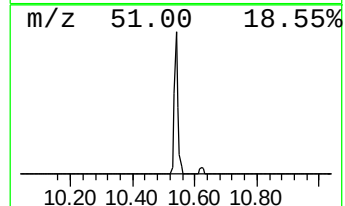
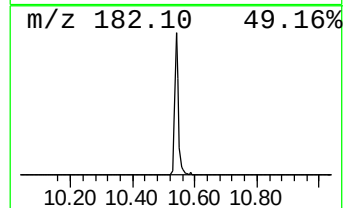
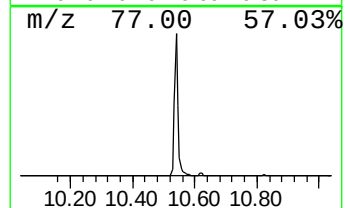
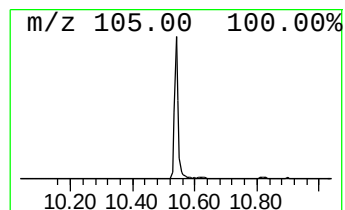
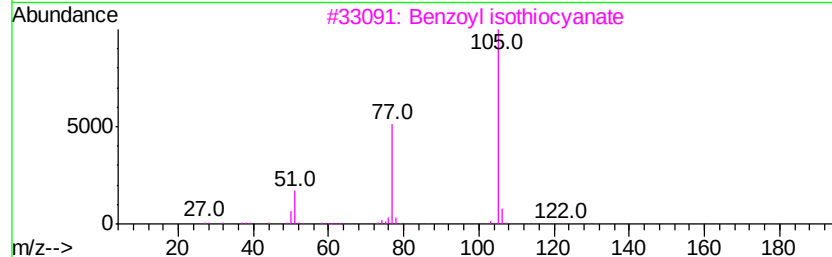
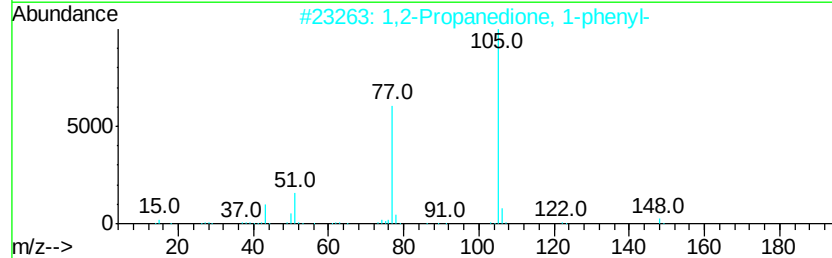
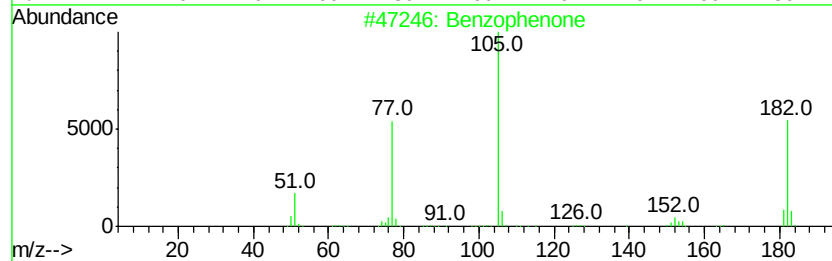
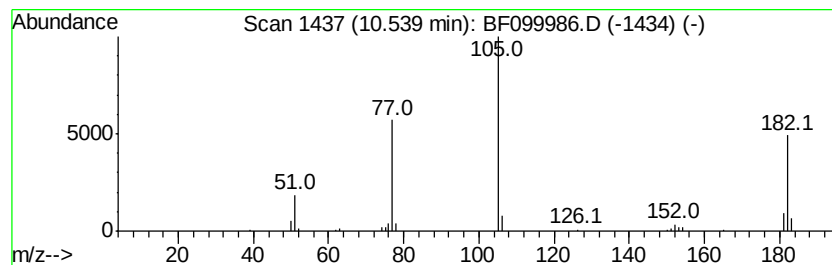
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Peak Number 4 Benzophenone Concentration Rank 3

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 10.54   | 3.87 ng | 127722                              | Acenaphthene-d10 | 9.82            |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | Benzophenone                        | 182 C13H10O      | 000119-61-9 96  |
| 2       |         | 1,2-Propanedione, 1-phenyl-         | 148 C9H8O2       | 000579-07-7 43  |
| 3       |         | Benzoyl isothiocyanate              | 163 C8H5NOS      | 000532-55-8 43  |
| 4       |         | Benzoic acid, pentafluorophenyl ... | 288 C13H5F5O2    | 1000360-67-9 43 |
| 5       |         | Benzenebutanoic acid, .gamma.-oxo-  | 178 C10H10O3     | 002051-95-8 43  |



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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |      |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|------|--------|------|
|                      |       |         |       |          | #                     | RT   | Resp   | Conc |
| 2-Pentanone, 4-hy... | 5.00  | 3.6     | ng    | 76808    | 1                     | 6.79 | 428917 | 20.0 |
| unknown6.56          | 6.56  | 55.4    | ng    | 1187340  | 1                     | 6.79 | 428917 | 20.0 |
| Benzophenone         | 10.54 | 3.9     | ng    | 127722   | 3                     | 9.82 | 659728 | 20.0 |