

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101215\  
 Data File : BF082226.D  
 Acq On : 12 Oct 2015 18:11  
 Operator : UM/IZ  
 Sample : G3990-06  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PO-005

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 1 % 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-BF101015.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         | 4.994       | 248           | 250         | 263          | rBV      | 325402         | 373015        | 15.06%          | 2.355%        |
| 2         | 5.417       | 285           | 287         | 290          | rBV      | 705948         | 733028        | 29.59%          | 4.628%        |
| 3         | 6.457       | 376           | 378         | 381          | rBV      | 414956         | 454231        | 18.34%          | 2.868%        |
| 4         | 6.594       | 388           | 390         | 393          | rBV      | 1488669        | 1436003       | 57.96%          | 9.066%        |
| 5         | 6.834       | 409           | 411         | 413          | rBV      | 304186         | 351497        | 14.19%          | 2.219%        |
| 6         | 6.983       | 422           | 424         | 427          | rBV      | 1383412        | 1431408       | 57.78%          | 9.037%        |
| 7         | 7.406       | 458           | 461         | 463          | rBV      | 867825         | 1159277       | 46.79%          | 7.319%        |
| 8         | 8.114       | 521           | 523         | 526          | rBV      | 496520         | 485876        | 19.61%          | 3.068%        |
| 9         | 9.200       | 615           | 618         | 620          | rBV      | 2544692        | 2477369       | 100.00%         | 15.641%       |
| 10        | 9.589       | 650           | 652         | 654          | rBV      | 89717          | 86833         | 3.51%           | 0.548%        |
| 11        | 9.875       | 674           | 677         | 679          | rBV      | 687537         | 609242        | 24.59%          | 3.846%        |
| 12        | 10.298      | 711           | 714         | 716          | rBV      | 99226          | 98949         | 3.99%           | 0.625%        |
| 13        | 10.663      | 743           | 746         | 749          | rBV      | 1409770        | 1484783       | 59.93%          | 9.374%        |
| 14        | 10.823      | 758           | 760         | 762          | rVB      | 67775          | 63266         | 2.55%           | 0.399%        |
| 15        | 11.360      | 805           | 807         | 809          | rBV      | 698647         | 624778        | 25.22%          | 3.945%        |
| 16        | 11.886      | 851           | 853         | 855          | rBV      | 46061          | 57121         | 2.31%           | 0.361%        |
| 17        | 11.955      | 855           | 859         | 861          | rVB      | 132003         | 184111        | 7.43%           | 1.162%        |
| 18        | 12.675      | 919           | 922         | 929          | rBV3     | 27852          | 50082         | 2.02%           | 0.316%        |
| 19        | 12.949      | 943           | 946         | 949          | rBV      | 2359851        | 2431129       | 98.13%          | 15.349%       |
| 20        | 13.841      | 1022          | 1024        | 1030         | rBV2     | 29891          | 42995         | 1.74%           | 0.271%        |
| 21        | 14.001      | 1035          | 1038        | 1044         | rBV      | 678514         | 683778        | 27.60%          | 4.317%        |
| 22        | 15.429      | 1160          | 1163        | 1169         | rBV      | 314880         | 520343        | 21.00%          | 3.285%        |

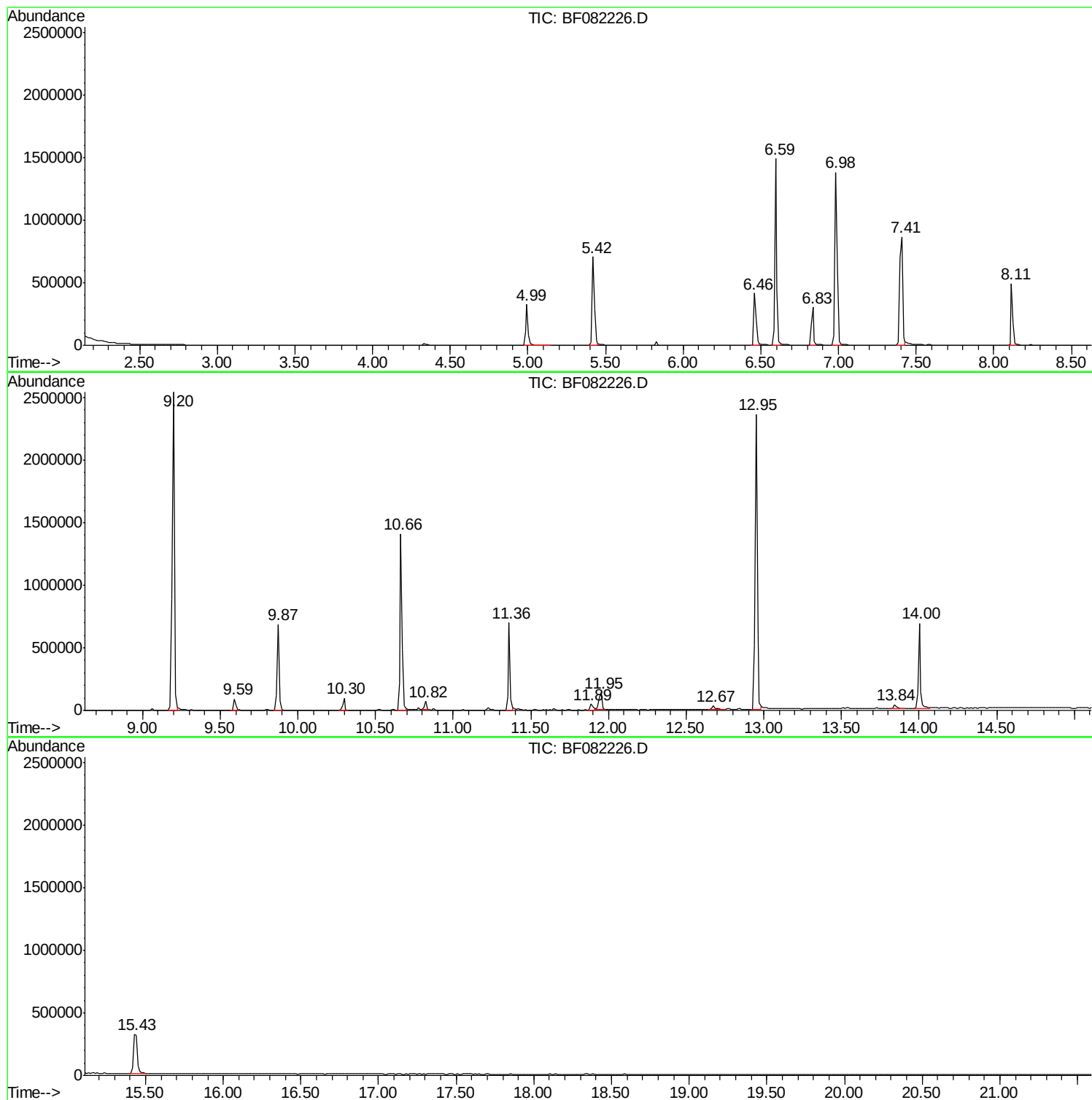
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BNA\_F  
ClientSampleId :  
PO-005

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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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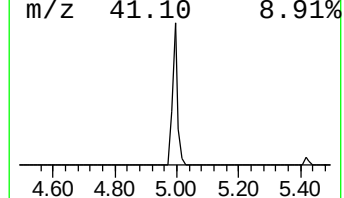
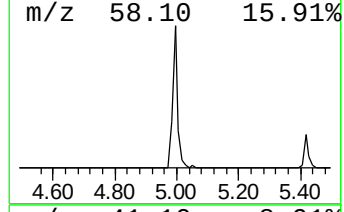
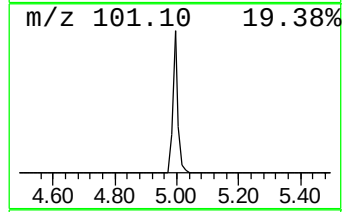
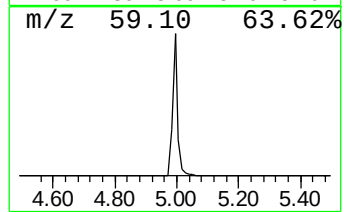
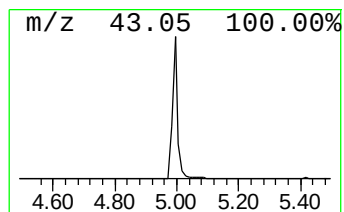
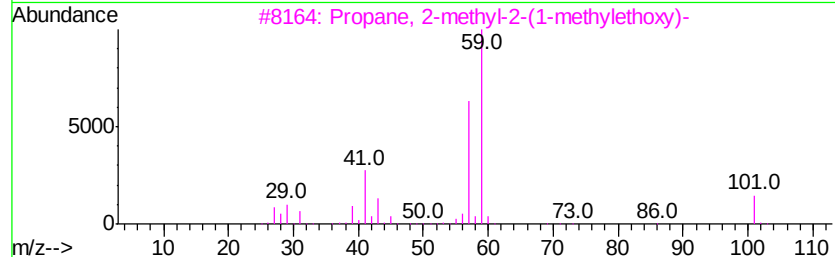
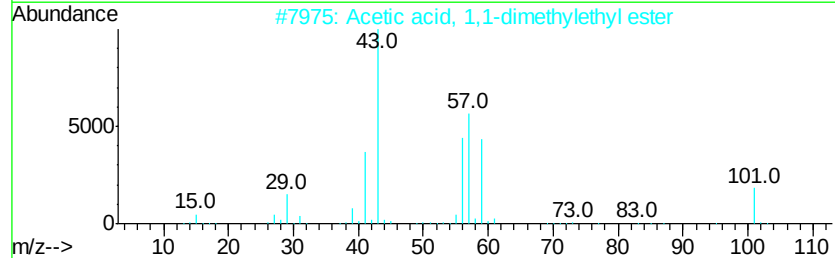
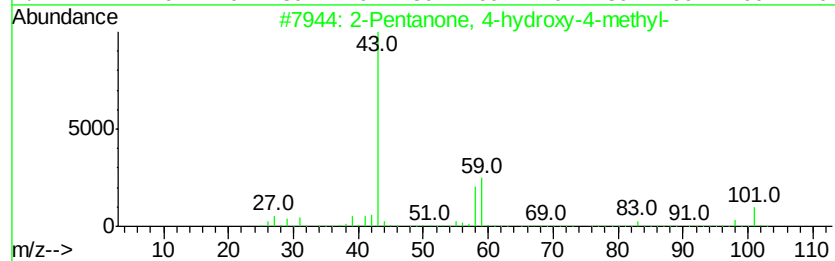
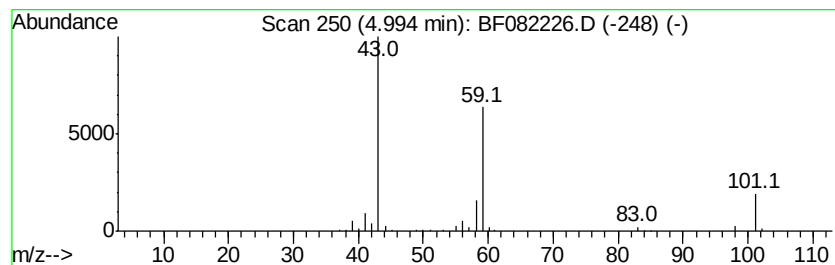
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.99 | 21.22 ng | 373015 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 50   |
| 2    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 36   |
| 4    |    | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 28   |
| 5    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |



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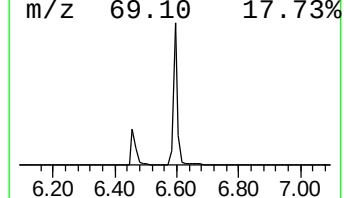
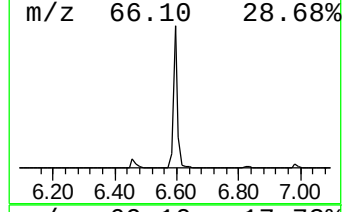
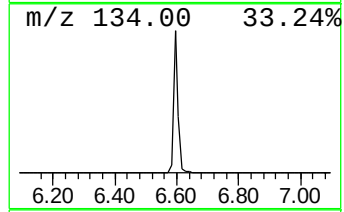
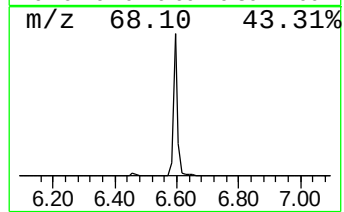
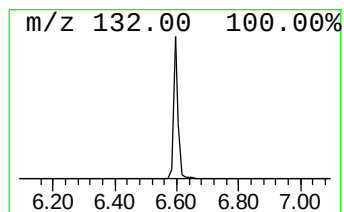
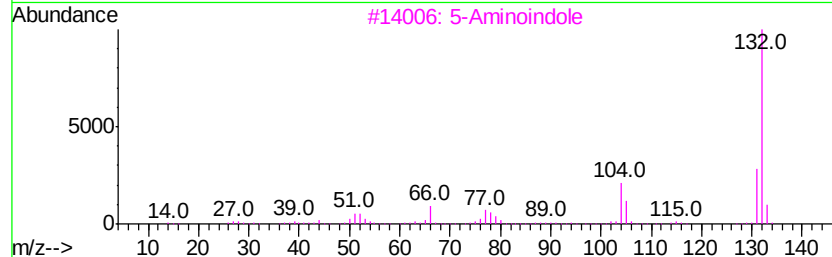
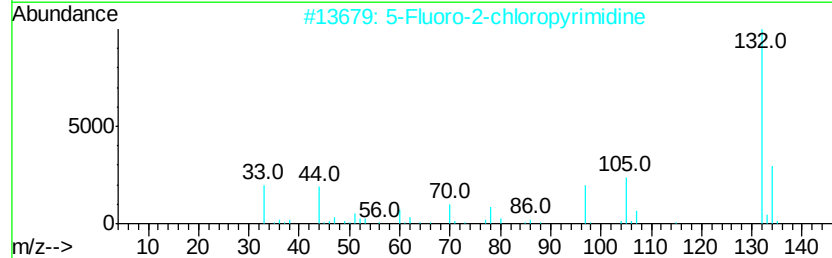
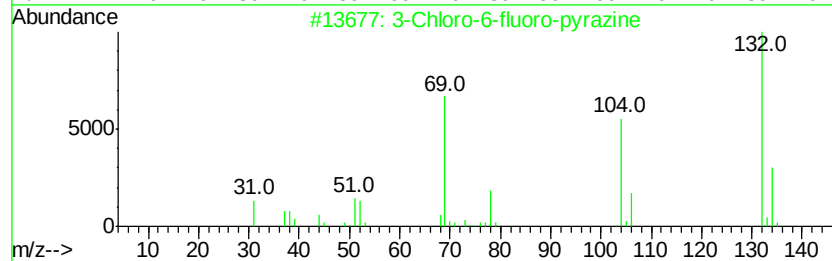
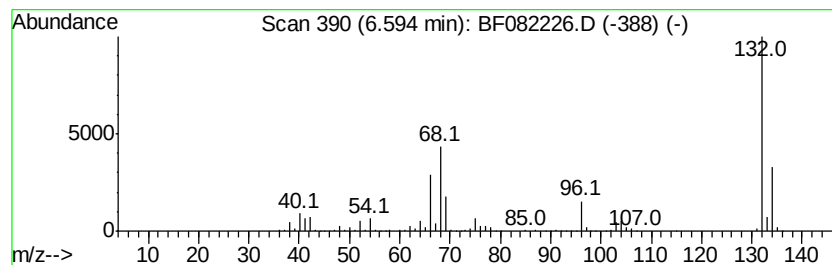
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TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 3-Chloro-6-fluoro-pyrazine Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.59 | 81.71 ng | 1436000 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    |   | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 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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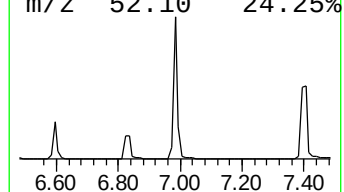
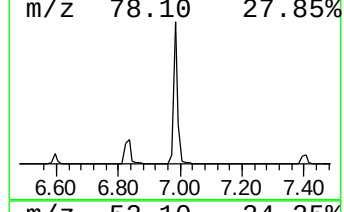
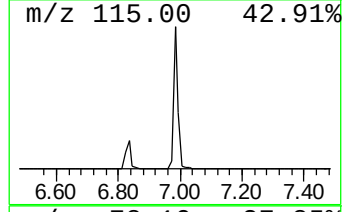
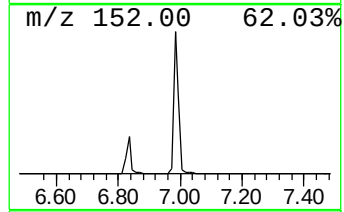
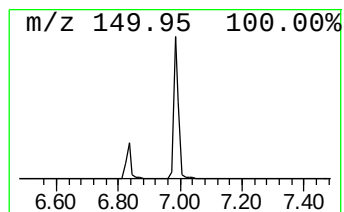
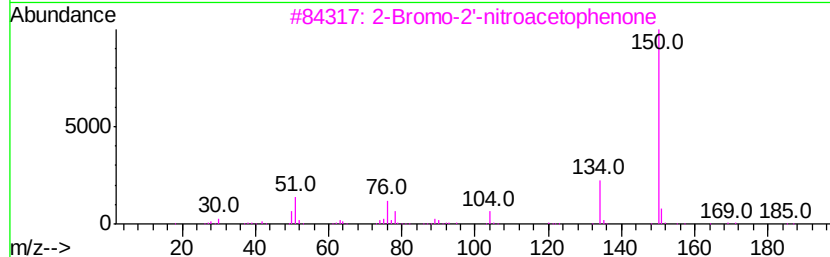
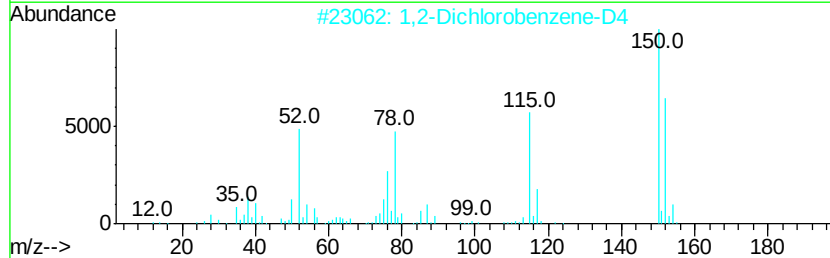
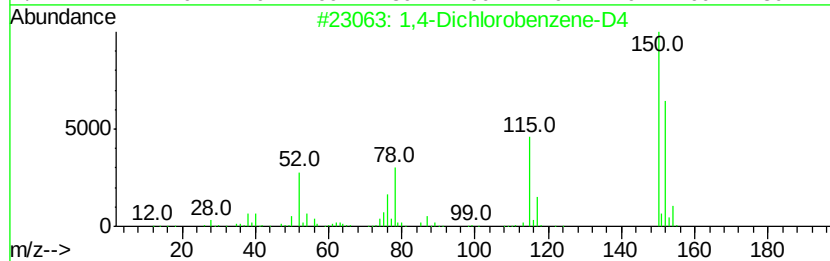
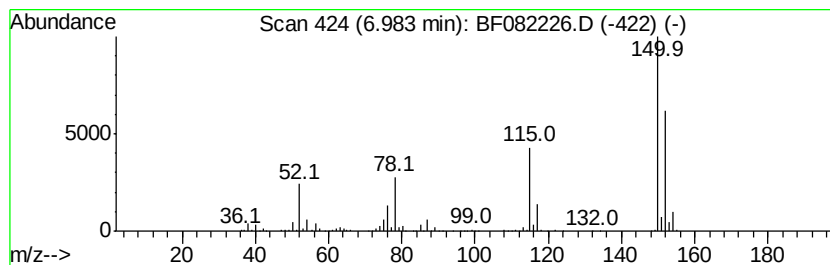
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 1,4-Dichlorobenzene-D4 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.98 | 81.45 ng | 1431410 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm       | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------------|-------------|------|
| 1    | 5  | 1,4-Dichlorobenzene-D4              | 150 | C6D4Cl2       | 003855-82-1 | 94   |
| 2    |    | 1,2-Dichlorobenzene-D4              | 150 | C6D4Cl2       | 002199-69-1 | 91   |
| 3    |    | 2-Bromo-2'-nitroacetophenone        | 243 | C8H6BrN03     | 006851-99-6 | 12   |
| 4    |    | 3-(4-Nitrobenzoyl)-2-thioxo-4-th... | 431 | C17H9N3O7S2   | 329218-74-8 | 10   |
| 5    |    | 2-Cyclohexene-1,4-dione, 5,6-dic... | 370 | C15H12Cl2N2O5 | 324051-56-1 | 10   |



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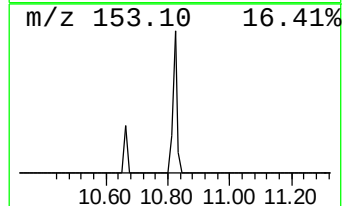
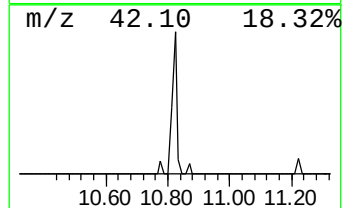
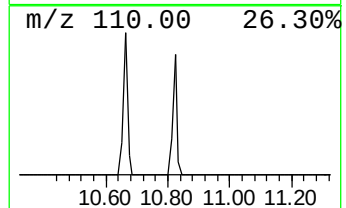
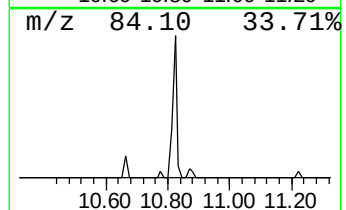
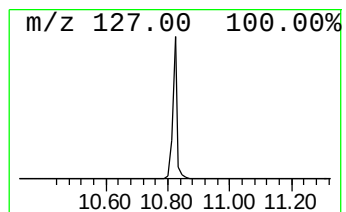
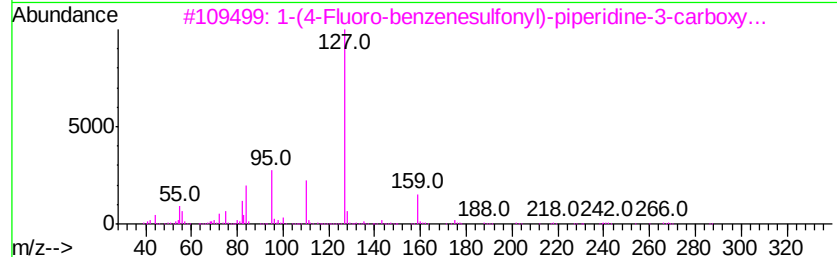
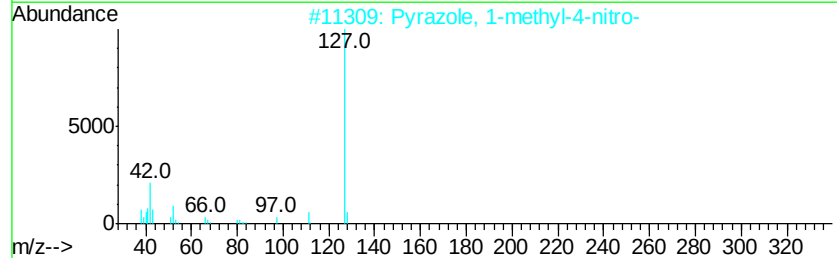
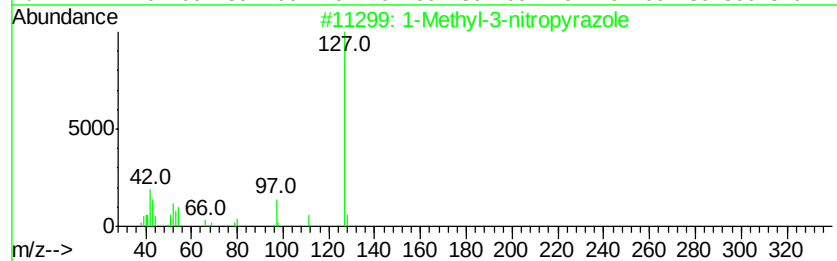
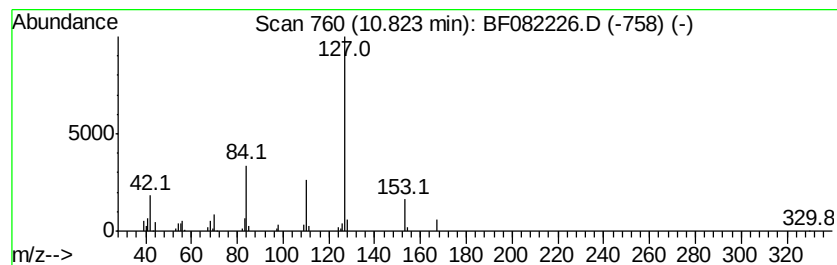
Instrument :  
BNA\_F  
ClientSampleId :  
PO-005

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\*\*\*\*\*  
Peak Number 4 1-Methyl-3-nitropyrazole Concentration Rank 5

| R.T.      | EstConc                              | Area  | Relative to ISTD | R.T.         |      |
|-----------|--------------------------------------|-------|------------------|--------------|------|
| 10.82     | 2.03 ng                              | 63266 | Phenanthrene-d10 | 11.36        |      |
| Hit# of 5 | Tentative ID                         | MW    | MolForm          | CAS#         | Qual |
| 1         | 1-Methyl-3-nitropyrazole             | 127   | C4H5N3O2         | 054210-32-1  | 46   |
| 2         | Pyrazole, 1-methyl-4-nitro-          | 127   | C4H5N3O2         | 003994-50-1  | 43   |
| 3         | 1-(4-Fluoro-benzenesulfonyl)-pip...  | 286   | C12H15FN2O3S     | 1000295-13-0 | 42   |
| 4         | Isopropylphosphonic acid, fluoroa... | 238   | C11H24F02P       | 1000273-54-8 | 40   |
| 5         | 1,3-Dimethylbutyl isopropylphosp...  | 210   | C9H20F02P        | 1000298-25-6 | 40   |



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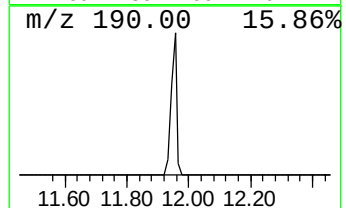
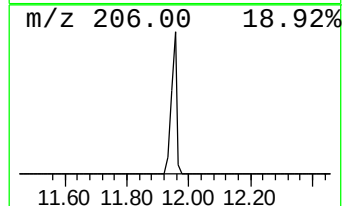
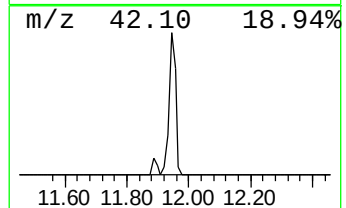
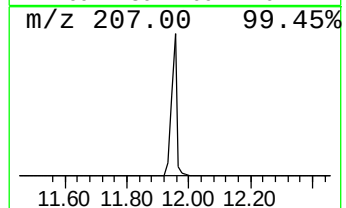
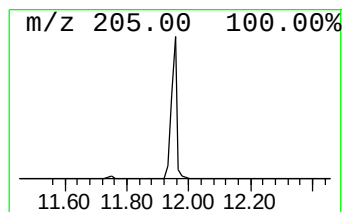
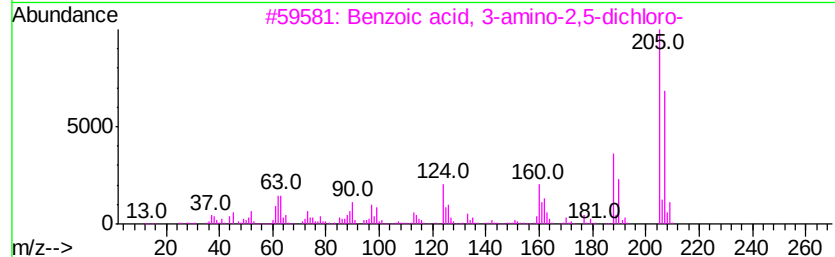
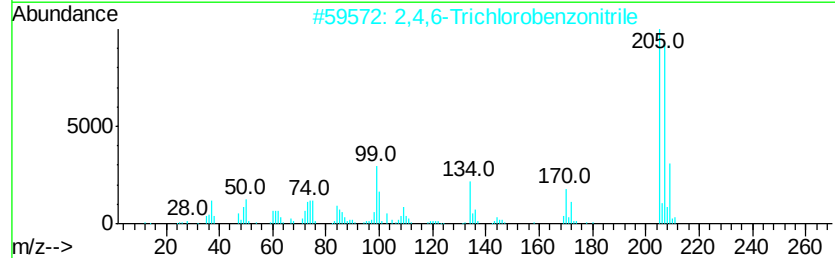
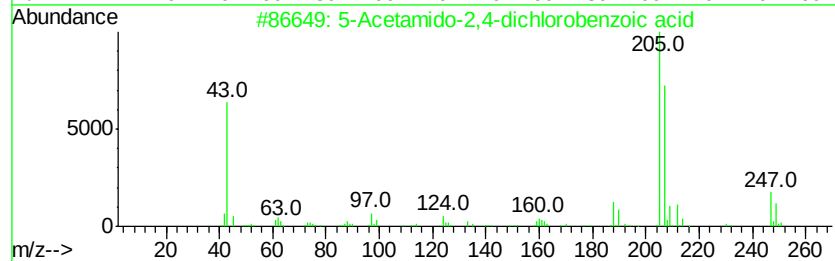
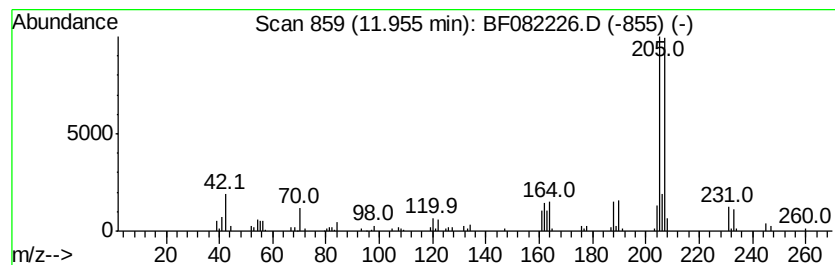
Instrument :  
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ClientSampleId :  
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Peak Number 5 5-Acetamido-2,4-dichloroben... Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD                    |     | R.T.       |             |      |
|-------|---------|--------|-------------------------------------|-----|------------|-------------|------|
| 11.95 | 5.89 ng | 184111 | Phenanthrene-d10                    |     | 11.36      |             |      |
| Hit#  | of      | 5      | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
| 1     |         |        | 5-Acetamido-2,4-dichlorobenzoic ... | 247 | C9H7Cl2NO3 | 050602-49-8 | 64   |
| 2     |         |        | 2,4,6-Trichlorobenzonitrile         | 205 | C7H2Cl3N   | 006575-05-9 | 59   |
| 3     |         |        | Benzoic acid, 3-amino-2,5-dichloro- | 205 | C7H5Cl2NO2 | 000133-90-4 | 50   |
| 4     |         |        | 5-Chloro-2,3-dihydrofuro(2,3-b)q... | 205 | C11H8ClNO  | 064124-92-1 | 38   |
| 5     |         |        | 7-Chloro-2,3-dihydrofuro(2,3-b)q... | 205 | C11H8ClNO  | 064124-91-0 | 37   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101215\  
Data File : BF082226.D  
Acq On : 12 Oct 2015 18:11  
Operator : UM/IZ  
Sample : G3990-06  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PO-005

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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 |
| 2-Pentanone, 4-hy... | 4.99  | 21.2    | ng    | 373015   | 1                     | 6.83  | 351497 | 20.0 |
| 3-Chloro-6-fluoro... | 6.59  | 81.7    | ng    | 1436000  | 1                     | 6.83  | 351497 | 20.0 |
| 1,4-Dichlorobenze... | 6.98  | 81.4    | ng    | 1431410  | 1                     | 6.83  | 351497 | 20.0 |
| 1-Methyl-3-nitrop... | 10.82 | 2.0     | ng    | 63266    | 4                     | 11.36 | 624778 | 20.0 |
| 5-Acetamido-2,4-d... | 11.95 | 5.9     | ng    | 184111   | 4                     | 11.36 | 624778 | 20.0 |