

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091415\  
 Data File : BF081619.D  
 Acq On : 14 Sep 2015 21:53  
 Operator : UM/IZ  
 Sample : G3597-13 5X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 GP-13(0-5)

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-BF090115.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         | 1.463       | 9             | 11          | 16           | rBV      | 22103          | 55162         | 1.30%           | 0.586%        |
| 2         | 1.543       | 16            | 18          | 70           | rVB      | 1188626        | 4256481       | 100.00%         | 45.225%       |
| 3         | 4.480       | 273           | 275         | 290          | rBV      | 58058          | 127290        | 2.99%           | 1.352%        |
| 4         | 5.383       | 352           | 354         | 362          | rBV      | 100133         | 205252        | 4.82%           | 2.181%        |
| 5         | 7.646       | 550           | 552         | 558          | rBV      | 107152         | 219055        | 5.15%           | 2.327%        |
| 6         | 7.749       | 558           | 561         | 570          | rVB      | 139950         | 245934        | 5.78%           | 2.613%        |
| 7         | 8.229       | 600           | 603         | 609          | rBB      | 205252         | 319262        | 7.50%           | 3.392%        |
| 8         | 8.549       | 629           | 631         | 637          | rBB      | 95425          | 167575        | 3.94%           | 1.780%        |
| 9         | 9.520       | 713           | 716         | 726          | rBB      | 85810          | 141406        | 3.32%           | 1.502%        |
| 10        | 11.132      | 854           | 857         | 861          | rBV      | 268145         | 413278        | 9.71%           | 4.391%        |
| 11        | 13.772      | 1085          | 1088        | 1091         | rBV      | 170635         | 247377        | 5.81%           | 2.628%        |
| 12        | 15.270      | 1215          | 1219        | 1223         | rBV2     | 252161         | 410639        | 9.65%           | 4.363%        |
| 13        | 17.167      | 1382          | 1385        | 1390         | rBB      | 63343          | 105483        | 2.48%           | 1.121%        |
| 14        | 18.779      | 1522          | 1526        | 1528         | rBV      | 209627         | 361668        | 8.50%           | 3.843%        |
| 15        | 18.824      | 1528          | 1530        | 1534         | rVB      | 74697          | 112030        | 2.63%           | 1.190%        |
| 16        | 21.442      | 1756          | 1759        | 1762         | rBV      | 181514         | 220024        | 5.17%           | 2.338%        |
| 17        | 21.796      | 1786          | 1790        | 1793         | rBV      | 161614         | 257045        | 6.04%           | 2.731%        |
| 18        | 22.116      | 1816          | 1818        | 1820         | rBV      | 167404         | 176167        | 4.14%           | 1.872%        |
| 19        | 23.396      | 1925          | 1930        | 1931         | rBV2     | 338681         | 479783        | 11.27%          | 5.098%        |
| 20        | 23.419      | 1931          | 1932        | 1935         | rVB      | 116852         | 126001        | 2.96%           | 1.339%        |
| 21        | 23.842      | 1964          | 1969        | 1971         | rBV2     | 35099          | 83677         | 1.97%           | 0.889%        |
| 22        | 24.493      | 2023          | 2026        | 2030         | rBV      | 130591         | 228264        | 5.36%           | 2.425%        |
| 23        | 24.733      | 2045          | 2047        | 2049         | rBV      | 70134          | 73993         | 1.74%           | 0.786%        |
| 24        | 24.779      | 2049          | 2051        | 2054         | rBV      | 74502          | 94407         | 2.22%           | 1.003%        |
| 25        | 24.836      | 2054          | 2056        | 2058         | rBV      | 161453         | 218937        | 5.14%           | 2.326%        |
| 26        | 25.888      | 2145          | 2148        | 2152         | rBV      | 35670          | 65607         | 1.54%           | 0.697%        |

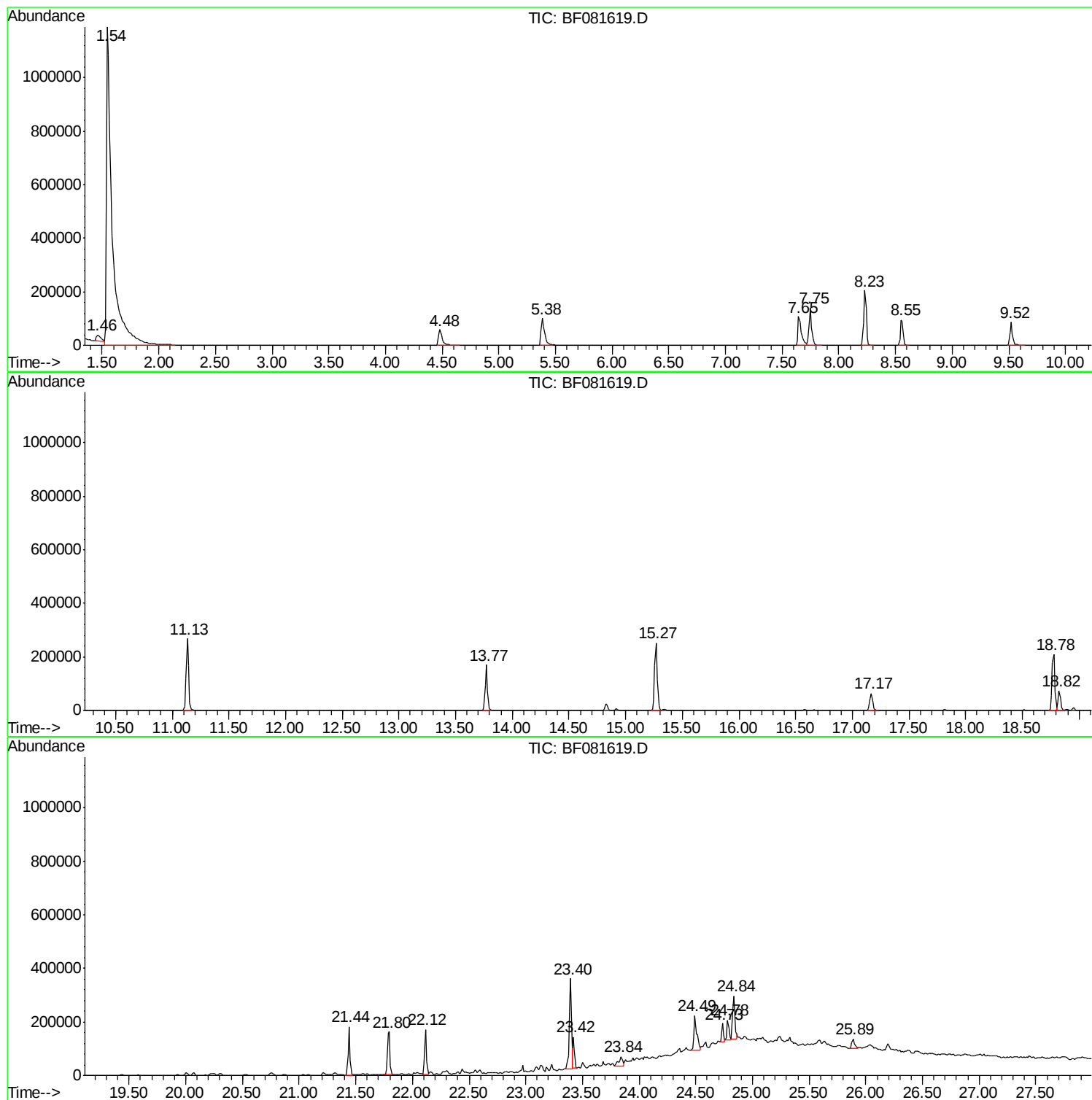
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ClientSampleId :  
GP-13(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF090115.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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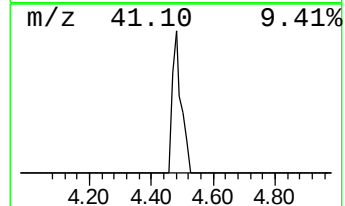
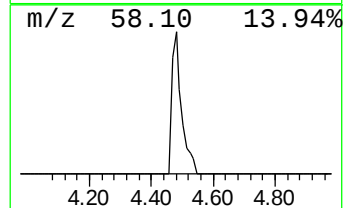
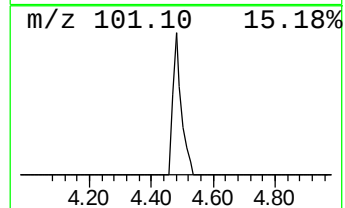
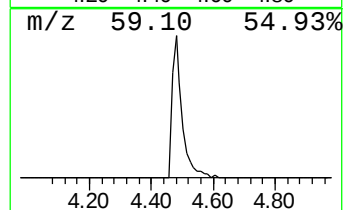
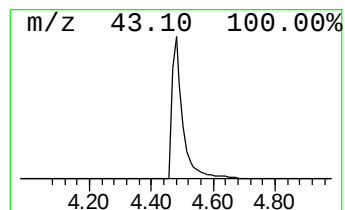
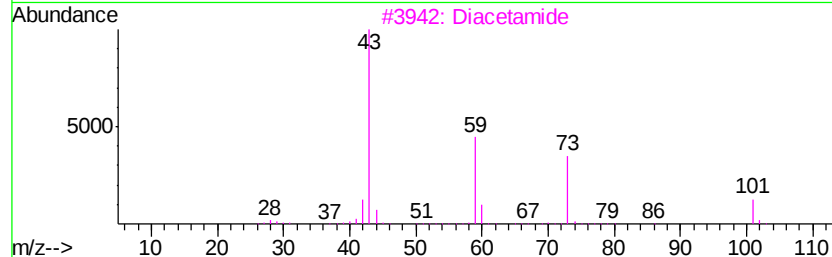
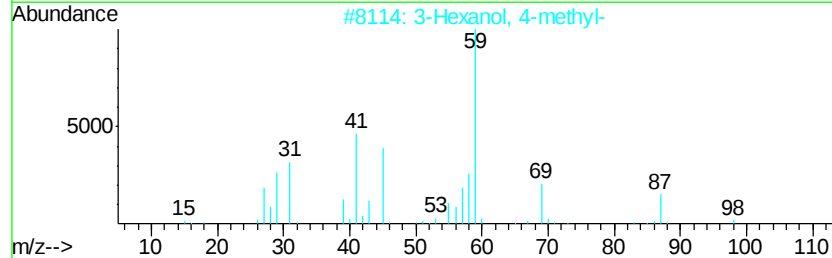
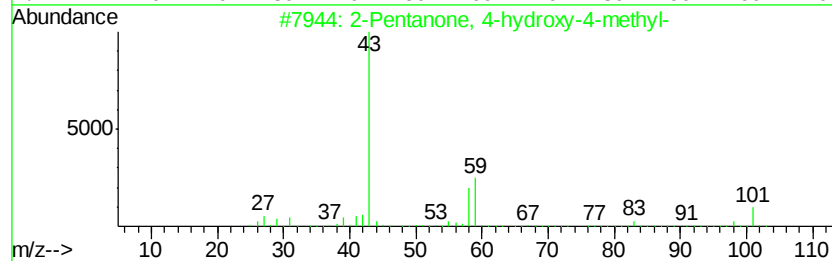
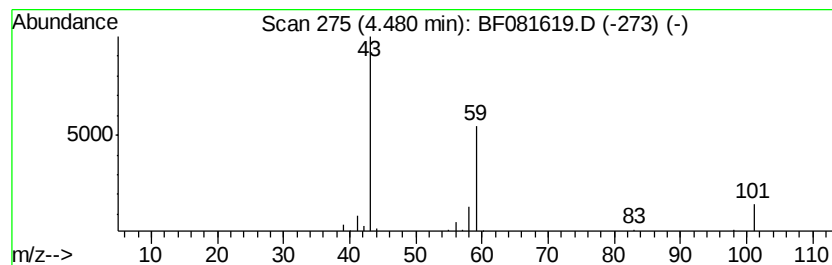
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.48 | 7.97 ng | 127290 | 1,4-Dichlorobenzene-d4 | 8.23 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 64   |
| 2    |      | 3-Hexanol, 4-methyl-                | 116 | C7H16O  | 000615-29-2 | 33   |
| 3    |      | Diacetamide                         | 101 | C4H7NO2 | 000625-77-4 | 9    |
| 4    |      | 1-Propen-2-ol, acetate              | 100 | C5H8O2  | 000108-22-5 | 9    |
| 5    |      | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O  | 017348-59-3 | 9    |



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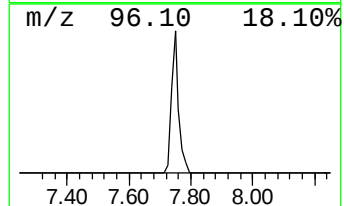
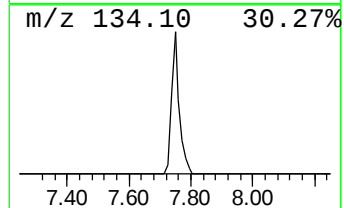
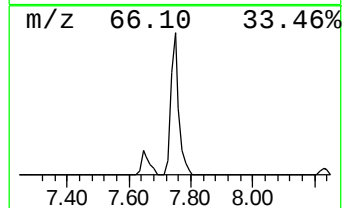
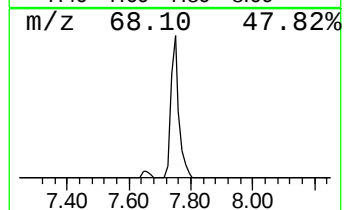
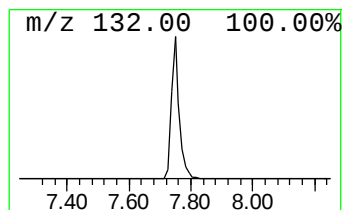
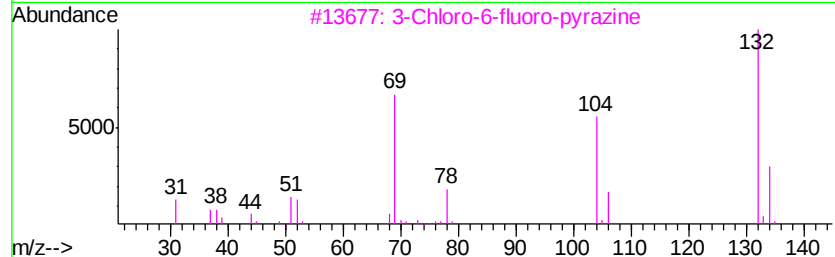
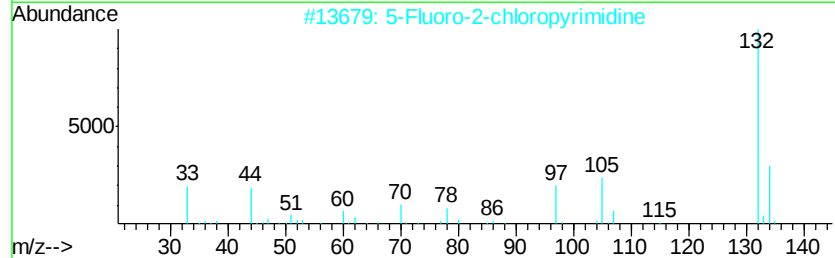
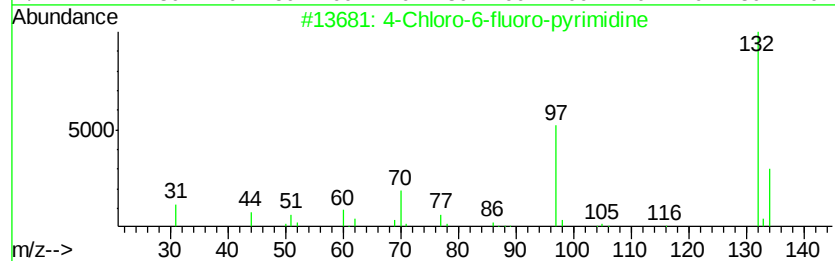
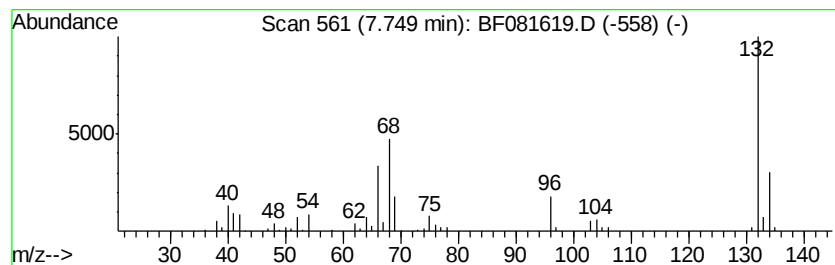
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Peak Number 4 unknown7.75 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.75 | 15.41 ng | 245934 | 1,4-Dichlorobenzene-d4 | 8.23 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#         | Qual |
|------|----|------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine       | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | 5-Fluoro-2-chloropyrimidine        | 132 | C4H2ClFN2 | 062802-42-0  | 27   |
| 3    |    | 3-Chloro-6-fluoro-pyrazine         | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 4    |    | Benzene, 1,3,5-trifluoro-          | 132 | C6H3F3    | 000372-38-3  | 14   |
| 5    |    | Pyrazolo(2,3-a)pyridine, 7-methyl- | 132 | C8H8N2    | 034760-58-2  | 12   |



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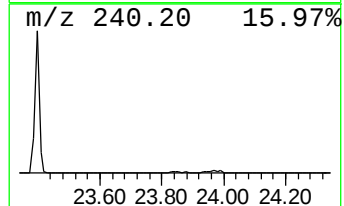
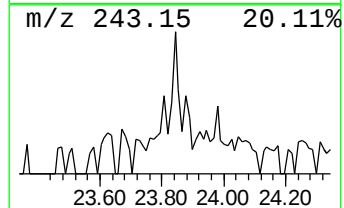
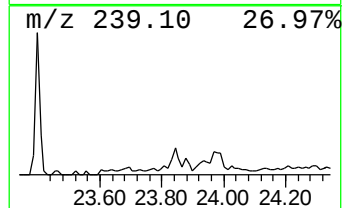
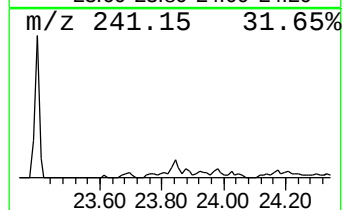
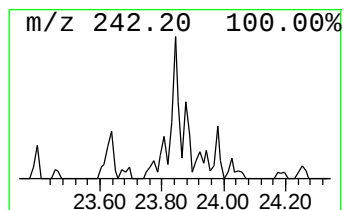
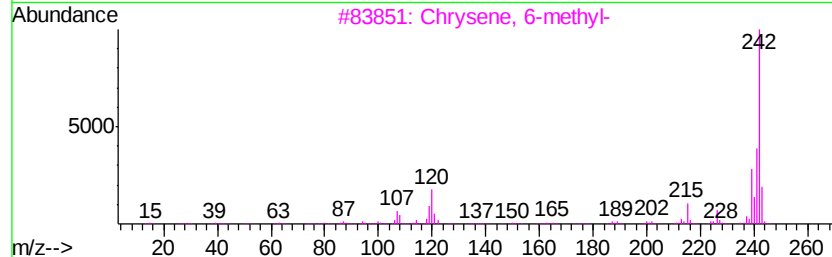
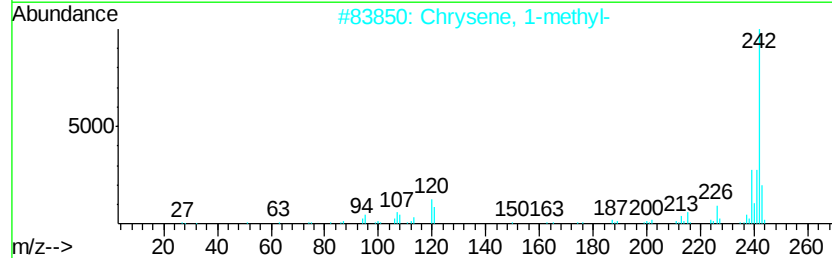
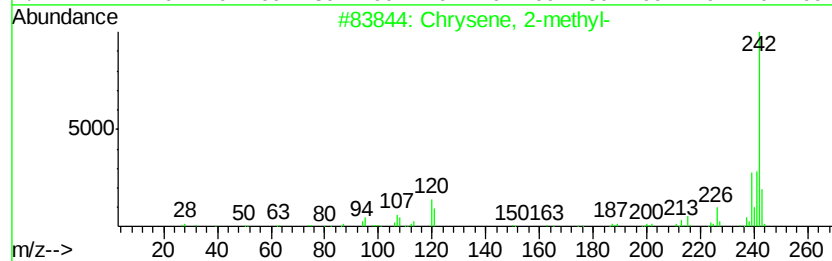
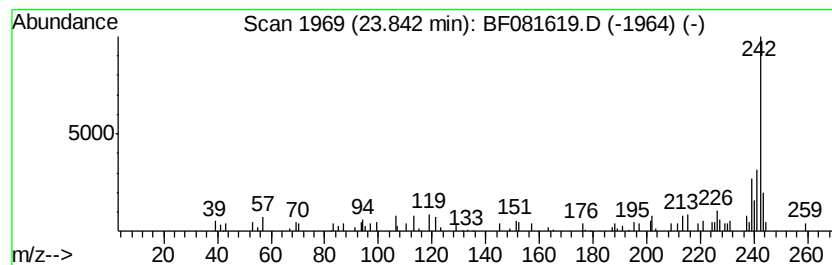
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BNA\_F  
ClientSampleId :  
GP-13(0-5)

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Peak Number 6 Chrysene, 2-methyl- Concentration Rank 5

| R.T.      | EstConc                      | Area  | Relative to ISTD |             | R.T.  |
|-----------|------------------------------|-------|------------------|-------------|-------|
| 23.84     | 3.49 ng                      | 83677 | Chrysene-d12     |             | 23.40 |
| Hit# of 5 | Tentative ID                 | MW    | MolForm          | CAS#        | Qual  |
| 1         | Chrysene, 2-methyl-          | 242   | C19H14           | 003351-32-4 | 81    |
| 2         | Chrysene, 1-methyl-          | 242   | C19H14           | 003351-28-8 | 81    |
| 3         | Chrysene, 6-methyl-          | 242   | C19H14           | 003697-24-3 | 76    |
| 4         | Benz[a]anthracene, 8-methyl- | 242   | C19H14           | 002381-31-9 | 74    |
| 5         | Triphenylene, 2-methyl-      | 242   | C19H14           | 001705-84-6 | 74    |



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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.48  | 8.0     | ng    | 127290   | 1                     | 8.23  | 319262 | 20.0 |
| unknown7.75          | 7.75  | 15.4    | ng    | 245934   | 1                     | 8.23  | 319262 | 20.0 |
| Chrysene, 2-methyl-  | 23.84 | 3.5     | ng    | 83677    | 5                     | 23.40 | 479783 | 20.0 |